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(54) Title: QUINUCLIDINE-SUBSTITUTED HETEROARYL MOIETIES FOR TREATMENT OF DISEASE

(57) Abstract: The invention provides compound of Formula (I): These compounds may be in the form of pharmaceutical salts or compositions, and racemic mixtures or pure enantiomers thereof. The compounds of Formula (I) are useful in pharmaceuticals in which  $\alpha 7$  is known to be involved.



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## QUINUCLIDINE-SUBSTITUTED HETEROARYL MOIETIES FOR TREATMENT OF DISEASE

### FIELD OF INVENTION

5           Nicotinic acetylcholine receptors (nAChRs) play a large role in central nervous system (CNS) activity. Particularly, they are known to be involved in cognition, learning, mood, emotion, and neuroprotection. There are several types of nicotinic acetylcholine receptors, and each one appears to have a different role in regulating CNS function. Nicotine affects all such receptors, and has a variety of activities.

10   Unfortunately, not all of the activities are desirable. In fact, one of the least desirable properties of nicotine is its addictive nature and the low ratio between efficacy and safety. The present invention relates to molecules that have a greater effect upon the  $\alpha 7$  nAChRs as compared to other closely related members of this large ligand-gated receptor family. Thus, the invention provides compounds that are active drug

15   molecules with fewer side effects.

### BACKGROUND OF THE INVENTION

          US Patent 5,977,144 discloses compositions for benzylidene- and cinnamylidene-anabaseines and methods for using these compositions for treating

20   conditions associated with defects or malfunctioning of nicotinic subtypes brain receptors. These compositions target the  $\alpha 7$  receptor subtype with little or no activation of the  $\alpha 4\beta 2$  or other receptor subtypes.

          US Patent 5,837,489 discloses human neuronal nicotinic acetylcholine receptor and cells transformed with same DNA and mRNA encoding subunits.

25           US Patent 5,712,270 discloses a group of 2-aroylaminothiazole derivatives which bind to and stimulate central muscarinic acetylcholine receptors and are useful agents for treating symptoms of cognitive disorders, specifically the impaired memory associated with a decrease in the neurotransmitter, acetylcholine. Some of the compounds of this invention also bind to 5HT<sub>1A</sub> receptors and dopamine D<sub>2</sub> receptors,

30   making them useful as antipsychotic agents.

          US Patent 5,624,941 discloses pyrazole derivatives useful in pharmaceuticals in which cannabis is known to be involved.

US Patent 5,561,149 discloses the use of a mono or bicyclic carbocyclic, or heterocyclic carboxylic, acid ester or amide or an imidazolyl carbazol in the manufacture of a medicament suitable for the treatment of stress-related psychiatric disorders, for increasing vigilance, for the treatment of rhinitis or serotonin-induced disorders and/or coadministration with another active agent to increase the bioavailability thereof, or for nasal administration.

US Patent 5,510,478 discloses a group of 2-arylaminothiazole derivatives which bind to and stimulate central muscarinic acetylcholine receptors and are useful agents for treating symptoms of cognitive disorders, specifically the impaired memory associated with a decrease in the neurotransmitter, acetylcholine. Some of the compounds of this invention also bind to 5HT<sub>1A</sub> receptors and dopamine D<sub>2</sub> receptors, making them useful as antipsychotic agents.

US Patent 5,364,863 discloses bicyclic carboxylic esters and amides, their pharmaceutical formulations, and a method for their use in treating migraine, emesis, gastrointestinal disorders, schizophrenia, or anxiety in mammals.

US Patent 5,342,845 discloses indole derivatives and drugs effective as gastrointestinal motor activity regulator, antimigraine, antipsychotic or antianxiety drugs.

US Patent 5,273,972 discloses novel 2-substituted-3-quinuclidinyl arylcarboxamides and arylthiocarboxamides and corresponding arylcarboxylates which have utility as therapeutic agents which exhibit gastric prokinetic, antiemetic, anxiolytic and 5-HT (serotonin) antagonist effects in warm blooded animals.

US Patent 5,246,942 discloses certain dibenzofurancarboxamides and their use as 5-HT<sub>3</sub> antagonists having unique CNS, anti-emetic and gastric prokinetic activity void of any significant D<sub>2</sub> receptor binding properties.

US Patent 5,237,066 discloses enantiomers of absolute configuration S of amide derivatives of 3-aminoquinuclidine, the process for preparing them and their use as medicinal products having activity in respect of gastric movements and antiemetic activity.

US Patent 5,236,931 discloses novel 3-quinuclidinyl benzamides and benzoates which have utility as therapeutical agents which exhibit anxiolytic, antipsychotic, cognition improvement, antiemetic and gastric prokinetic effects in warm blooded animals.

US Patent 5,217,975 discloses azabicyclic compounds for treating dementia.

US Patent 5,206,246 discloses anxiolytic-R-N-(1-azabicyclo[2.2.2]oct-3-yl) benzamides and thiobenzamides, their N-oxides and pharmaceutically acceptable salts thereof. A preferred compound is R-(+)-4-amino-N-(1-azabicyclo[2.2.2]oct-3-yl)-5-  
5 chloro-2-methoxybenzamide.

US Patent 5,183,822 discloses new heterocyclic compounds (3,4-annulated benzimidazole-2(1H)-ones) having an antagonistic activity on 5-hydroxytryptamine (5-HT) receptors.

US Patent 5,175,173 discloses carboxamides useful as antiemetic or  
10 antipsychotic agents.

US Patent 5,106,843 discloses heterocyclic compounds useful as 5-HT<sub>3</sub> antagonists.

US Patent 5,070,095 discloses novel 1-(azabicyclo[2.2.2]oct-3- or -4-yl)benzamides substituted on the benzene ring with the basic substituted  
15 aminomethyleneamino group which have been found to be useful in treating emesis, including emesis due to chemical and radiation anticancer therapy, anxiety, and impaired gastric emptying.

US Patent 5,057,519 discloses 5-HT<sub>3</sub> antagonists as being useful in reducing opiate tolerance.

US Patent 5,039,680 discloses 5-HT<sub>3</sub> antagonists in preventing or reducing  
20 dependency on dependency-inducing agents.

US Patent 5,025,022 discloses a method of treating or preventing schizophrenia and/or psychosis using S-N-(1-azabicyclo[2.2.2]oct-3-yl)benzamides and thiobenzamides, their N-oxides and pharmaceutically acceptable salts thereof. A  
25 preferred compound is S-(-)-4-amino-N-(1-azabicyclo[2.2.2]oct-3-yl)-5-chloro-2-methoxybenzamide.

US Patent 5,017,580 discloses memory enhancing R-N-(1-azabicyclo[2.2.2.]oct-3-yl)benzamides and thiobenzamides, their N-oxides and pharmaceutically acceptable salts thereof. A preferred compound is R-(+)-4-amino-  
30 N-(1-azabicyclo[2.2.2]oct-3-yl)-5-chloro-2-methoxybenzamide.

US Patent 4,988,691 discloses isoxazole-containing compounds exhibiting anti-serotonin activity.

US Patent 4,921,982 discloses 5-halo-2,3-dihydro-2,2-dimethylbenzofuran-7-carboxylic acids which are useful as intermediates for 5-HT<sub>3</sub> antagonists.

US Patent 4,835,162 discloses agonists and antagonists to nicotine as smoking deterrents.

5        US Patent 4,822,795 discloses pharmaceutically useful esters and amides having 5-HT<sub>3</sub> antagonist activity.

US Patent 4,803,199 discloses pharmaceutically useful heterocyclic acid esters and amides or alkylene bridged piperidines as serotonin M antagonists.

10        US Patent 4,798,829 discloses 1-azabicyclo[3.2.2]nonane derivatives having gastric motility enhancing activity and/or anti-emetic activity and/or 5-HT receptor antagonist activity.

US Patent 4,789,673 discloses dicarboxylic, heterocyclic and substituted benzoic acid alkylene-bridged piperidyl amides and esters as being serotonin M antagonists.

15        US Patent 4,721,720 discloses a method of treating emesis, anxiety and/or irritable bowel syndrome.

US Patent 4,657,911 discloses 3-amino quinuclidine derivatives and the application thereof as accelerators of gastro-intestinal motor function and as medicament potentiators.

20        US Patent 4,605,652 discloses a method of enhancing memory or correcting memory deficiency with arylamido (and arylthioamido)-azabicycloalkanes, and the pharmaceutically acceptable acid addition salts, hydrates and alcoholates thereof.

US Patent 3,702,324 discloses 3,4,5-trimethoxybenzamides of substituted anilines and of alkylpiperidines which exert a specific effect on the central nervous system and a somewhat lesser effect on muscle function, and thus have utility as tranquilizers.

25        WO 01/36417 A1 discloses novel N-azabicyclo-amide derivatives and use in therapy, especially in the treatment of prophylaxis of psychotic disorders and intellectual impairment disorders.

30        WO 00/73431 A2 discloses two binding assays to directly measure the affinity and selectivity of compounds at the  $\alpha 7$  nAChR and the 5-HT<sub>3</sub>R. The combined use of these functional and binding assays may be used to identify compounds that are selective agonists of the  $\alpha 7$  nAChR.

WO 92/15579 discloses multicyclic tertiary amine polyaromatic squalene synthase inhibitors and method of treatment for lowering serum cholesterol levels using the compounds.

WO 92/11259 discloses azabicyclic amides or esters of halogenated benzoic acids having 5-HT<sub>3</sub> receptor antagonist activity.

WO 90/14347 A as abstracted in chemical abstract 1991:143,158 discloses N-quinuclidinyl-indolecarboxamide derivatives as being antiemetics.

EP 512 350 A2 discloses 3-(indolyl-2-carboxamido) quinuclidines useful for treating diseases characterized by an excess or enhanced sensitivity to serotonin, e.g., psychosis, nausea, vomiting, dementia or other cognitive diseases, migraine, diabetes. The compound may be used to control anxiety, aggression, depression, and pain. The compounds are disclosed as serotonin 5-HT<sub>3</sub> antagonists.

DE 3810552 A1 discloses esters and amides of indolyl-, benzo[b]thiophenyl-, benzo[b]furancarboxylic acids or 4-amino-2 methoxy-benzoic acids with N-heterocyclic or N-heterobicyclic alcohols or amines. The compounds disclosed have activity against pain especially migraine, as an anti-arrhythmic for gastrointestinal disturbances, stomach disturbances, gastritis ulcer, gall bladder, spastic colon, Crohn's disease, ulcerative colitis, carcinoid syndrome, diarrhea of various types. The compounds are also disclosed as speeding stomach emptying, controlling gastro duodenal and gastro esophageal reflux, disturbances of esophageal motility, hiatal hernia, cardiac insufficiency, hypotonic stomach, paralytic ileus, manic depressive psychosis and other psychoses. The compounds are also disclosed as useful for stress related diseases, senility, and enhancement of nasal absorption of other agents, e.g., in the treatment of emesis.

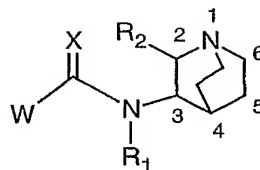
FR 2 625 678 discloses N-(quinuclidin-3-yl)-benzamides and thiobenzamides useful as diet-control agents.

In *Bioorg. & Med.Chem. Lett.* 11 (2001) 319-321, the 5-HT<sub>3</sub> antagonist tropisetron (ICS 205-930) is discussed as a potent and selective  $\alpha 7$  Nicotinic receptor partial agonist.

In *Behavioral Brain Res.*, 113 (2000) 169-181, it is discussed that the brain  $\alpha 7$  nicotinic receptor may be an important therapeutic target for the treatment of Alzheimer's disease using DMXBA which is known as GTS-21.

## SUMMARY OF THE INVENTION

A compound of Formula I:



Formula I

5 or pharmaceutically acceptable salts thereof wherein

$R_1$  is selected from -H, alkyl, cycloalkyl, halogenated alkyl, or aryl;

Alkyl is both straight and branched-chain moieties having from 1-6 carbon atoms;

Halogenated alkyl is an alkyl moiety having from 1-6 carbon atoms and having  
10 1 to (2n+1) substituent(s) independently selected from -F, -Cl, -Br, or -I where n is the maximum number of carbon atoms in the moiety;

Cycloalkyl is a cyclic alkyl moiety having from 3-6 carbon atoms;

Aryl is phenyl, substituted phenyl, naphthyl, or substituted naphthyl;

Substituted phenyl is a phenyl either having 1-4 substituents independently  
15 selected from -F, -Cl, -Br, or -I, or having 1 substituent selected from  $-R_{12}$  and 0-3 substituents independently selected from -F, -Cl, -Br, or -I;

Substituted naphthyl is a naphthalene moiety either having 1-4 substituents independently selected from -F, -Cl, -Br, or -I, or having 1 substituent selected from  
20  $-R_{12}$  and 0-3 substituents independently selected from -F, -Cl, -Br, or -I, where the substitution can be independently on either the same ring or different rings of said naphthalene moiety;

$R_2$  is -H, alkyl, halogenated alkyl, substituted alkyl, cycloalkyl, benzyl, substituted benzyl, or aryl;

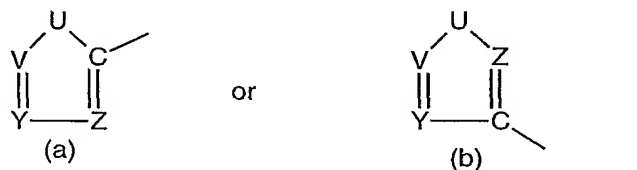
Substituted alkyl is an alkyl moiety having from 1-6 carbon atoms and having  
25 0-3 substituents independently selected from -F, -Cl, -Br, or -I, and further having 1 substituent selected from  $-R_7$ ,  $-R_9$ ,  $-OR_{10}$ ,  $-SR_{10}$ ,  $-NR_{10}R_{10}$ ,  $-C(O)R_{10}$ ,  $-NO_2$ ,  $-C(O)NR_{10}R_{10}$ ,  $-CN$ ,  $-NR_{10}C(O)R_{10}$ ,  $-S(O)_2NR_{10}R_{10}$ ,  $-NR_{10}S(O)_2R_{10}$ , phenyl, or substituted phenyl;

Substituted benzyl is a benzyl either having 1-4 substituents independently  
30 selected from -F, -Cl, -Br, or -I, or having 1 substituent selected from  $-R_{12}$  and 0-3

substituents independently selected from -F, -Cl, -Br, or -I, provided that all substitution is on the phenyl ring of the benzyl;

X is O or S;

W is a cyclic heteroaromatic moiety where the heteroatoms can be from 1-3 atoms selected from oxygen, sulfur, or nitrogen of the following structures:



wherein U is -O-, -S-, or -N(R<sub>3</sub>)-;

V and Y are independently selected from =N-, or =C(R<sub>5</sub>)-;

Z is =N-, or =CH-, provided that when both V and Y are =C(R<sub>5</sub>)- and Z is =CH-, only one =C(R<sub>5</sub>)- can be =CH-, and further provided that when U is -O-, Y is =C(R<sub>5</sub>)- and Z is =C(R<sub>5</sub>)-, V cannot be =N-,

R<sub>3</sub> is -H, alkyl, cycloalkyl, heterocycloalkyl, halogenated alkyl, halogenated cycloalkyl, halogenated heterocycloalkyl, substituted alkyl, limited substituted alkyl, substituted cycloalkyl, substituted heterocycloalkyl, or aryl, and provided that when W is (b) and Z is =N- and U is N(R<sub>3</sub>), R<sub>3</sub> cannot be phenyl or substituted phenyl;

Limited substituted alkyl is a substituted alkyl having from 1-6 carbon atoms and having 0-3 substituents independently selected from -F, -Cl, -Br, or -I, and further having 1 substituent on either only the ω carbon and selected from -OR<sub>11</sub>, -SR<sub>11</sub>, -NR<sub>11</sub>R<sub>11</sub>, -C(O)R<sub>11</sub>, -NO<sub>2</sub>, -C(O)NR<sub>11</sub>R<sub>11</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>11</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, or -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, or on any carbon with sufficient valency but not on the ω carbon and selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -NO<sub>2</sub>, -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl;

Alkenyl is straight and branched-chain moieties having from 2-6 carbon atoms and having at least one carbon-carbon double bond;

Halogenated alkenyl is an unsaturated alkenyl moiety having from 2-6 carbon atoms and having 1 to (2n-1) substituent(s) independently selected from -F, -Cl, -Br, or -I where n is the maximum number of carbon atoms in the moiety;

Substituted alkenyl is an unsaturated alkenyl moiety having from 2-6 carbon atoms and having 0-3 substituents independently selected from -F, or -Cl, and further having 1 substituent selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>,

-C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl;

Alkynyl is straight and branched-chain moieties having from 2-6 carbon atoms and having at least one carbon-carbon triple bond;

5 Halogenated alkynyl is an unsaturated alkynyl moiety having from 3-6 carbon atoms and having 1 to (2n-3) substituent(s) independently selected from -F, -Cl, -Br, or -I where n is the maximum number of carbon atoms in the moiety;

Substituted alkynyl is an unsaturated alkynyl moiety having from 3-6 carbon atoms and having 0-3 substituents independently selected from -F, or -Cl, and further  
10 having 1 substituent selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -CN, -C(O)NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl;

Halogenated cycloalkyl is a cyclic moiety having from 3-6 carbon atoms and having 1-4 substituents independently selected from -F, or -Cl;

15 Substituted cycloalkyl is a cyclic moiety having from 3-6 carbon atoms and having 0-3 substituents independently selected from -F, or -Cl, and further having 1 substituent selected from -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, -NO<sub>2</sub>, phenyl, or substituted phenyl;

Heterocycloalkyl is a cyclic moiety having 4-7 atoms with 1-2 atoms within  
20 the ring being -S-, -N(R<sub>3</sub>)-, or -O-;

Halogenated heterocycloalkyl is a cyclic moiety having from 4-7 atoms with 1-2 atoms within the ring being -S-, -N(R<sub>3</sub>)-, or -O-, and having 1-4 substituents independently selected from -F, or -Cl;

Substituted heterocycloalkyl is a cyclic moiety having from 4-7 atoms with 1-2  
25 atoms within the ring being -S-, -N(R<sub>3</sub>)-, or -O- and having 0-3 substituents independently selected from -F, or -Cl, and further having 1 substituent selected from -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -NO<sub>2</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl;

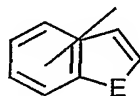
R<sub>5</sub> is independently selected from the group consisting of -H, alkyl, alkenyl, alkyne, cycloalkyl, heterocycloalkyl, halogenated alkyl, halogenated alkenyl, halogenated alkyne, halogenated cycloalkyl, halogenated heterocycloalkyl, substituted alkyl, substituted alkenyl, substituted alkyne, substituted cycloalkyl, substituted heterocycloalkyl, limited substituted alkyl, limited substituted alkenyl,

limited substituted alkynyl, aryl,  $-OR_8$ ,  $-OR_{14}$ ,  $-SR_8$ ,  $-SR_{14}$ ,  $-F$ ,  $-Cl$ ,  $-Br$ ,  $-I$ ,  $-NR_8R_8$ ,  $-NR_{14}R_{14}$ ,  $-C(O)R_8$ ,  $-C(O)R_{14}$ ,  $-C(O)NR_8R_8$ ,  $-C(O)NR_{14}R_{14}$ ,  $-CN$ ,  $-NR_8C(O)R_{11}$ ,  $-S(O)_2NR_8R_8$ ,  $-OS(O)_2R_{11}$ ,  $-S(O)_2R_8$ ,  $-S(O)_2R_{14}$ ,  $-NR_8S(O)_2R_8$ ,  $-N(H)C(O)N(H)R_8$ ,  $-NO_2$ ,  $-R_7$ , and  $-R_9$ ;

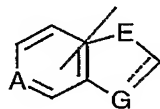
5 Limited substituted alkenyl is a substituted alkenyl having from 1-6 carbon atoms and having 0-3 substituents independently selected from  $-F$ ,  $-Cl$ ,  $-Br$ , or  $-I$ , and further having 1 substituent on either only the  $\omega$  carbon and selected from  $-OR_{11}$ ,  $-SR_{11}$ ,  $-NR_{11}R_{11}$ ,  $-C(O)R_{11}$ ,  $-NO_2$ ,  $-C(O)NR_{11}R_{11}$ ,  $-CN$ ,  $-NR_{10}C(O)R_{11}$ ,  $-S(O)_2NR_{10}R_{10}$ , or  $-NR_{10}S(O)_2R_{10}$ , or on any carbon with sufficient valency but not on  
10 the  $\omega$  carbon and selected from  $-R_7$ ,  $-R_9$ ,  $-OR_{10}$ ,  $-SR_{10}$ ,  $-NR_{10}R_{10}$ ,  $-C(O)R_{10}$ ,  $-NO_2$ ,  $-C(O)NR_{10}R_{10}$ ,  $-CN$ ,  $-NR_{10}C(O)R_{10}$ ,  $-S(O)_2NR_{10}R_{10}$ ,  $-NR_{10}S(O)_2R_{10}$ , phenyl, or substituted phenyl;

Limited substituted alkynyl is a substituted alkynyl having from 1-6 carbon atoms and having 0-3 substituents independently selected from  $-F$ ,  $-Cl$ ,  $-Br$ , or  $-I$ , and  
15 further having 1 substituent on either only the  $\omega$  carbon and selected from  $-OR_{11}$ ,  $-SR_{11}$ ,  $-NR_{11}R_{11}$ ,  $-C(O)R_{11}$ ,  $-NO_2$ ,  $-C(O)NR_{11}R_{11}$ ,  $-CN$ ,  $-NR_{10}C(O)R_{11}$ ,  $-S(O)_2NR_{10}R_{10}$ , or  $-NR_{10}S(O)_2R_{10}$ , or on any carbon with sufficient valency but not on the  $\omega$  carbon and selected from  $-R_7$ ,  $-R_9$ ,  $-OR_{10}$ ,  $-SR_{10}$ ,  $-NR_{10}R_{10}$ ,  $-C(O)R_{10}$ ,  $-NO_2$ ,  $-C(O)NR_{10}R_{10}$ ,  $-CN$ ,  $-NR_{10}C(O)R_{10}$ ,  $-S(O)_2NR_{10}R_{10}$ ,  $-NR_{10}S(O)_2R_{10}$ , phenyl, or  
20 substituted phenyl;

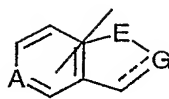
$R_7$  is 5-membered heteroaromatic mono-cyclic moieties containing within the ring 1-3 heteroatoms independently selected from the group consisting of  $-O-$ ,  $=N-$ ,  $-N(R_3)-$ , and  $-S-$ , and having 0-1 substituent selected from  $-R_{12}$  and 0-3 substituents independently selected from  $-F$ ,  $-Cl$ ,  $-Br$ , or  $-I$ , or  $R_7$  is a 9-membered fused-ring  
25 moiety having a 6-membered ring fused to a 5-membered ring and having the formula



wherein E is O, S, or  $NR_3$ ,



wherein E and G are independently selected from  $CR_{18}$ , O, S, or  $NR_3$ , and A is  $CR_{18}$   
30 or N, or



wherein E and G are independently selected from  $\text{CR}_{18}$ , O, S, or  $\text{NR}_3$ , and A is  $\text{CR}_{18}$  or N, each 9-membered fused-ring moiety having 0-1 substituent selected from  $-\text{R}_{12}$  and 0-3 substituent(s) independently selected from -F, -Cl, -Br, or -I, and having a bond directly or indirectly attached to the core molecule where valency allows in either the 6-membered or the 5-membered ring of the fused-ring moiety;

Each  $\text{R}_8$  is independently selected from -H, alkyl, halogenated alkyl, substituted alkyl, cycloalkyl, halogenated cycloalkyl, substituted cycloalkyl, heterocycloalkyl, halogenated heterocycloalkyl, substituted heterocycloalkyl,  $-\text{R}_7$ ,  $-\text{R}_9$ , phenyl, or substituted phenyl;

$\text{R}_9$  is 6-membered heteroaromatic mono-cyclic moieties containing within the ring 1-3 heteroatoms selected from =N- and having 0-1 substituent selected from  $-\text{R}_{12}$  and 0-3 substituent(s) independently selected from -F, -Cl, -Br, or -I, or 10-membered heteroaromatic bi-cyclic moieties containing within one or both rings 1-3 heteroatoms selected from =N-, including, but not limited to, quinolinyl or isoquinolinyl, each 10-membered fused-ring moiety having 0-1 substituent selected from  $-\text{R}_{12}$  and 0-3 substituent(s) independently selected from -F, -Cl, -Br, or -I and having a bond directly or indirectly attached to the core molecule where valency allows;

Each  $\text{R}_{10}$  is independently selected from -H, alkyl, cycloalkyl, heterocycloalkyl, alkyl substituted with 1 substituent selected from  $\text{R}_{13}$ , cycloalkyl substituted with 1 substituent selected from  $\text{R}_{13}$ , heterocycloalkyl substituted with 1 substituent selected from  $\text{R}_{13}$ , halogenated alkyl, halogenated cycloalkyl, halogenated heterocycloalkyl, phenyl, substituted phenyl,  $-\text{R}_7$ , or  $-\text{R}_9$ ;

Each  $\text{R}_{11}$  is independently selected from -H, alkyl, cycloalkyl, heterocycloalkyl, halogenated alkyl, halogenated cycloalkyl, or halogenated heterocycloalkyl;

$\text{R}_{12}$  is selected from  $-\text{OR}_{11}$ ,  $-\text{SR}_{11}$ , alkyl, cycloalkyl, heterocycloalkyl, halogenated alkyl, halogenated cycloalkyl, halogenated heterocycloalkyl, substituted alkyl, substituted cycloalkyl, substituted heterocycloalkyl,  $-\text{NR}_{11}\text{R}_{11}$ ,  $-\text{C(O)}\text{R}_{11}$ ,  $-\text{NO}_2$ ,  $-\text{C(O)}\text{NR}_{11}\text{R}_{11}$ , -CN,  $-\text{NR}_{11}\text{C(O)}\text{R}_{11}$ ,  $-\text{S(O)}_2\text{NR}_{11}\text{R}_{11}$ , or  $-\text{NR}_{11}\text{S(O)}_2\text{R}_{11}$ ;

$\text{R}_{13}$  is selected from  $-\text{OR}_{11}$ ,  $-\text{SR}_{11}$ ,  $-\text{NR}_{11}\text{R}_{11}$ ,  $-\text{C(O)}\text{R}_{11}$ ,  $-\text{C(O)}\text{NR}_{11}\text{R}_{11}$ , -CN,  $-\text{NR}_{11}\text{C(O)}\text{R}_{11}$ ,  $-\text{S(O)}_2\text{NR}_{11}\text{R}_{11}$ ,  $-\text{NR}_{11}\text{S(O)}_2\text{R}_{11}$ ,  $-\text{CF}_3$ , or  $-\text{NO}_2$ ;

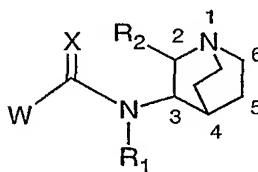
$R_{14}$  is independently selected from -H, alkyl, halogenated alkyl, limited substituted alkyl, cycloalkyl, halogenated cycloalkyl, substituted cycloalkyl, heterocycloalkyl, halogenated heterocycloalkyl, substituted heterocycloalkyl; and

- Each  $R_{18}$  is independently selected from -H, alkyl, cycloalkyl,  
 5 heterocycloalkyl, halogenated alkyl, halogenated cycloalkyl, halogenated heterocycloalkyl, substituted alkyl, substituted cycloalkyl, substituted heterocycloalkyl, -OR<sub>11</sub>, -SR<sub>11</sub>, -NR<sub>11</sub>R<sub>11</sub>, -C(O)R<sub>11</sub>, -NO<sub>2</sub>, -C(O)NR<sub>11</sub>R<sub>11</sub>, -CN, -NR<sub>11</sub>C(O)R<sub>11</sub>, -S(O)<sub>2</sub>NR<sub>11</sub>R<sub>11</sub>, or -NR<sub>11</sub>S(O)<sub>2</sub>R<sub>11</sub>, -F, -Cl, -Br, or -I, or a bond directly or indirectly attached to the core molecule, provided that there is only one said  
 10 bond to the core molecule within the 9-membered fused-ring moiety, further provided that the fused-ring moiety has 0-1 substituent selected from alkyl, cycloalkyl, heterocycloalkyl, halogenated alkyl, halogenated cycloalkyl, halogenated heterocycloalkyl, substituted alkyl, substituted cycloalkyl, substituted heterocycloalkyl, -OR<sub>11</sub>, -SR<sub>11</sub>, -NR<sub>11</sub>R<sub>11</sub>, -C(O)R<sub>11</sub>, -NO<sub>2</sub>, -C(O)NR<sub>11</sub>R<sub>11</sub>, -CN, -NR<sub>11</sub>C(O)R<sub>11</sub>, -S(O)<sub>2</sub>NR<sub>11</sub>R<sub>11</sub>, or -NR<sub>11</sub>S(O)<sub>2</sub>R<sub>11</sub>, and further provided that the fused-ring moiety has 0-3 substituent(s) selected from -F, -Cl, -Br, or -I.

Compounds of Formula I are useful to treat any one of or combination of schizophrenia, or psychosis.

## 20 DETAILED DESCRIPTION OF THE INVENTION

Surprisingly, we have found that compounds of Formula I:



Formula I

wherein  $R_1$  is selected from -H, alkyl, cycloalkyl, halogenated alkyl, or aryl;

- 25 Alkyl is both straight and branched-chain moieties having from 1-6 carbon atoms;

Halogenated alkyl is an alkyl moiety having from 1-6 carbon atoms and having 1 to (2n+1) substituent(s) independently selected from -F, -Cl, -Br, or -I where n is the maximum number of carbon atoms in the moiety;

- 30 Cycloalkyl is a cyclic alkyl moiety having from 3-6 carbon atoms;

Aryl is phenyl, substituted phenyl, naphthyl, or substituted naphthyl;

Substituted phenyl is a phenyl either having 1-4 substituents independently selected from -F, -Cl, -Br, or -I, or having 1 substituent selected from -R<sub>12</sub> and 0-3 substituents independently selected from -F, -Cl, -Br, or -I;

5 Substituted naphthyl is a naphthalene moiety either having 1-4 substituents independently selected from -F, -Cl, -Br, or -I, or having 1 substituent selected from -R<sub>12</sub> and 0-3 substituents independently selected from -F, -Cl, -Br, or -I, where the substitution can be independently on either the same ring or different rings of said naphthalene moiety;

10 R<sub>2</sub> is -H, alkyl, halogenated alkyl, substituted alkyl, cycloalkyl, benzyl, substituted benzyl, or aryl;

Substituted alkyl is an alkyl moiety having from 1-6 carbon atoms and having 0-3 substituents independently selected from -F, -Cl, -Br, or -I, and further having 1 substituent selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -NO<sub>2</sub>,

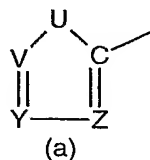
15 -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl;

Substituted benzyl is a benzyl either having 1-4 substituents independently selected from -F, -Cl, -Br, or -I, or having 1 substituent selected from -R<sub>12</sub> and 0-3 substituents independently selected from -F, -Cl, -Br, or -I, provided that all

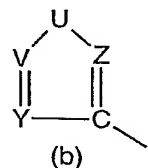
20 substitution is on the phenyl ring of the benzyl;

X is O or S;

W is a cyclic heteroaromatic moiety where the heteroatoms can be from 1-3 atoms selected from oxygen, sulfur, or nitrogen of the following structures:



or



25 wherein U is -O-, -S-, or -N(R<sub>3</sub>)-;

V and Y are independently selected from =N-, or =C(R<sub>5</sub>)-;

Z is =N-, or =CH-, provided that when both V and Y are =C(R<sub>5</sub>)- and Z is =CH-, only one =C(R<sub>5</sub>)- can be =CH-, and further provided that when U is -O-, Y is =C(R<sub>5</sub>)- and Z is =C(R<sub>5</sub>)-, V cannot be =N-,

$R_3$  is -H, alkyl, cycloalkyl, heterocycloalkyl, halogenated alkyl, halogenated cycloalkyl, halogenated heterocycloalkyl, substituted alkyl, limited substituted alkyl, substituted cycloalkyl, substituted heterocycloalkyl, or aryl, and provided that when W is (b) and Z is =N- and U is N( $R_3$ ),  $R_3$  cannot be phenyl or substituted phenyl;

5 Limited substituted alkyl is a substituted alkyl having from 1-6 carbon atoms and having 0-3 substituents independently selected from -F, -Cl, -Br, or -I, and further having 1 substituent on either only the  $\omega$  carbon and selected from -OR<sub>11</sub>, -SR<sub>11</sub>, -NR<sub>11</sub>R<sub>11</sub>, -C(O)R<sub>11</sub>, -NO<sub>2</sub>, -C(O)NR<sub>11</sub>R<sub>11</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>11</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, or -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, or on any carbon with sufficient valency but not on the  $\omega$  carbon and  
10 selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -NO<sub>2</sub>, -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl;

Alkenyl is straight and branched-chain moieties having from 2-6 carbon atoms and having at least one carbon-carbon double bond;

Halogenated alkenyl is an unsaturated alkenyl moiety having from 2-6 carbon  
15 atoms and having 1 to (2n-1) substituent(s) independently selected from -F, -Cl, -Br, or -I where n is the maximum number of carbon atoms in the moiety;

Substituted alkenyl is an unsaturated alkenyl moiety having from 2-6 carbon atoms and having 0-3 substituents independently selected from -F, or -Cl, and further having 1 substituent selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>,  
20 -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl;

Alkynyl is straight and branched-chain moieties having from 2-6 carbon atoms and having at least one carbon-carbon triple bond;

Halogenated alkynyl is an unsaturated alkynyl moiety having from 3-6 carbon  
25 atoms and having 1 to (2n-3) substituent(s) independently selected from -F, -Cl, -Br, or -I where n is the maximum number of carbon atoms in the moiety;

Substituted alkynyl is an unsaturated alkynyl moiety having from 3-6 carbon atoms and having 0-3 substituents independently selected from -F, or -Cl, and further having 1 substituent selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -CN,  
30 -C(O)NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl;

Halogenated cycloalkyl is a cyclic moiety having from 3-6 carbon atoms and having 1-4 substituents independently selected from -F, or -Cl;

Substituted cycloalkyl is a cyclic moiety having from 3-6 carbon atoms and having 0-3 substituents independently selected from -F, or -Cl, and further having 1 substituent selected from -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, -NO<sub>2</sub>, phenyl, or substituted phenyl;

5 Heterocycloalkyl is a cyclic moiety having 4-7 atoms with 1-2 atoms within the ring being -S-, -N(R<sub>3</sub>)-, or -O-;

Halogenated heterocycloalkyl is a cyclic moiety having from 4-7 atoms with 1-2 atoms within the ring being -S-, -N(R<sub>3</sub>)-, or -O-, and having 1-4 substituents independently selected from -F, or -Cl;

10 Substituted heterocycloalkyl is a cyclic moiety having from 4-7 atoms with 1-2 atoms within the ring being -S-, -N(R<sub>3</sub>)-, or -O- and having 0-3 substituents independently selected from -F, or -Cl, and further having 1 substituent selected from -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -NO<sub>2</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl;

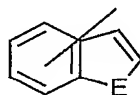
15 R<sub>5</sub> is independently selected from the group consisting of -H, alkyl, alkenyl, alkynyl, cycloalkyl, heterocycloalkyl, halogenated alkyl, halogenated alkenyl, halogenated alkynyl, halogenated cycloalkyl, halogenated heterocycloalkyl, substituted alkyl, substituted alkenyl, substituted alkynyl, substituted cycloalkyl, substituted heterocycloalkyl, limited substituted alkyl, limited substituted alkenyl, limited substituted alkynyl, aryl, -OR<sub>8</sub>, -OR<sub>14</sub>, -SR<sub>8</sub>, -SR<sub>14</sub>, -F, -Cl, -Br, -I, -NR<sub>8</sub>R<sub>8</sub>, -NR<sub>14</sub>R<sub>14</sub>, -C(O)R<sub>8</sub>, -C(O)R<sub>14</sub>, -C(O)NR<sub>8</sub>R<sub>8</sub>, -C(O)NR<sub>14</sub>R<sub>14</sub>, -CN, -NR<sub>8</sub>C(O)R<sub>11</sub>, -S(O)<sub>2</sub>NR<sub>8</sub>R<sub>8</sub>, -OS(O)<sub>2</sub>R<sub>11</sub>, -S(O)<sub>2</sub>R<sub>8</sub>, -S(O)<sub>2</sub>R<sub>14</sub>, -NR<sub>8</sub>S(O)<sub>2</sub>R<sub>8</sub>, -N(H)C(O)N(H)R<sub>8</sub>, -NO<sub>2</sub>, -R<sub>7</sub>, and -R<sub>9</sub>;

25 Limited substituted alkenyl is a substituted alkenyl having from 1-6 carbon atoms and having 0-3 substituents independently selected from -F, -Cl, -Br, or -I, and further having 1 substituent on either only the ω carbon and selected from -OR<sub>11</sub>, -SR<sub>11</sub>, -NR<sub>11</sub>R<sub>11</sub>, -C(O)R<sub>11</sub>, -NO<sub>2</sub>, -C(O)NR<sub>11</sub>R<sub>11</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>11</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, or -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, or on any carbon with sufficient valency but not on the ω carbon and selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -NO<sub>2</sub>, -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl;

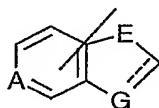
Limited substituted alkynyl is a substituted alkynyl having from 1-6 carbon atoms and having 0-3 substituents independently selected from -F, -Cl, -Br, or -I, and

further having 1 substituent on either only the  $\omega$  carbon and selected from  $-\text{OR}_{11}$ ,  $-\text{SR}_{11}$ ,  $-\text{NR}_{11}\text{R}_{11}$ ,  $-\text{C}(\text{O})\text{R}_{11}$ ,  $-\text{NO}_2$ ,  $-\text{C}(\text{O})\text{NR}_{11}\text{R}_{11}$ ,  $-\text{CN}$ ,  $-\text{NR}_{10}\text{C}(\text{O})\text{R}_{11}$ ,  $-\text{S}(\text{O})_2\text{NR}_{10}\text{R}_{10}$ , or  $-\text{NR}_{10}\text{S}(\text{O})_2\text{R}_{10}$ , or on any carbon with sufficient valency but not on the  $\omega$  carbon and selected from  $-\text{R}_7$ ,  $-\text{R}_9$ ,  $-\text{OR}_{10}$ ,  $-\text{SR}_{10}$ ,  $-\text{NR}_{10}\text{R}_{10}$ ,  $-\text{C}(\text{O})\text{R}_{10}$ ,  $-\text{NO}_2$ ,  
 5  $-\text{C}(\text{O})\text{NR}_{10}\text{R}_{10}$ ,  $-\text{CN}$ ,  $-\text{NR}_{10}\text{C}(\text{O})\text{R}_{10}$ ,  $-\text{S}(\text{O})_2\text{NR}_{10}\text{R}_{10}$ ,  $-\text{NR}_{10}\text{S}(\text{O})_2\text{R}_{10}$ , phenyl, or substituted phenyl;

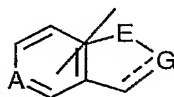
$\text{R}_7$  is 5-membered heteroaromatic mono-cyclic moieties containing within the ring 1-3 heteroatoms independently selected from the group consisting of  $-\text{O}-$ ,  $=\text{N}-$ ,  $-\text{N}(\text{R}_3)-$ , and  $-\text{S}-$ , and having 0-1 substituent selected from  $-\text{R}_{12}$  and 0-3 substituents  
 10 independently selected from  $-\text{F}$ ,  $-\text{Cl}$ ,  $-\text{Br}$ , or  $-\text{I}$ , or  $\text{R}_7$  is a 9-membered fused-ring moiety having a 6-membered ring fused to a 5-membered ring and having the formula



wherein E is O, S, or  $\text{NR}_3$ ,



15 wherein E and G are independently selected from  $\text{CR}_{18}$ , O, S, or  $\text{NR}_3$ , and A is  $\text{CR}_{18}$  or N, or



wherein E and G are independently selected from  $\text{CR}_{18}$ , O, S, or  $\text{NR}_3$ , and A is  $\text{CR}_{18}$  or N, each 9-membered fused-ring moiety having 0-1 substituent selected from  $-\text{R}_{12}$  and 0-3 substituent(s) independently selected from  $-\text{F}$ ,  $-\text{Cl}$ ,  $-\text{Br}$ , or  $-\text{I}$ , and having a  
 20 bond directly or indirectly attached to the core molecule where valency allows in either the 6-membered or the 5-membered ring of the fused-ring moiety;

Each  $\text{R}_8$  is independently selected from  $-\text{H}$ , alkyl, halogenated alkyl, substituted alkyl, cycloalkyl, halogenated cycloalkyl, substituted cycloalkyl,  
 25 heterocycloalkyl, halogenated heterocycloalkyl, substituted heterocycloalkyl,  $-\text{R}_7$ ,  $-\text{R}_9$ , phenyl, or substituted phenyl;

$\text{R}_9$  is 6-membered heteroaromatic mono-cyclic moieties containing within the ring 1-3 heteroatoms selected from  $=\text{N}-$  and having 0-1 substituent selected from  $-\text{R}_{12}$  and 0-3 substituent(s) independently selected from  $-\text{F}$ ,  $-\text{Cl}$ ,  $-\text{Br}$ , or  $-\text{I}$ , or 10-membered

heteroaromatic bi-cyclic moieties containing within one or both rings 1-3 heteroatoms selected from =N-, including, but not limited to, quinolinyl or isoquinolinyl, each 10-membered fused-ring moiety having 0-1 substituent selected from -R<sub>12</sub> and 0-3 substituent(s) independently selected from -F, -Cl, -Br, or -I and having a bond  
 5 directly or indirectly attached to the core molecule where valency allows;

Each R<sub>10</sub> is independently selected from -H, alkyl, cycloalkyl, heterocycloalkyl, alkyl substituted with 1 substituent selected from R<sub>13</sub>, cycloalkyl substituted with 1 substituent selected from R<sub>13</sub>, heterocycloalkyl substituted with 1 substituent selected from R<sub>13</sub>, halogenated alkyl, halogenated cycloalkyl, halogenated  
 10 heterocycloalkyl, phenyl, substituted phenyl, -R<sub>7</sub>, or -R<sub>9</sub>;

Each R<sub>11</sub> is independently selected from -H, alkyl, cycloalkyl, heterocycloalkyl, halogenated alkyl, halogenated cycloalkyl, or halogenated heterocycloalkyl;

R<sub>12</sub> is selected from -OR<sub>11</sub>, -SR<sub>11</sub>, alkyl, cycloalkyl, heterocycloalkyl, halogenated alkyl, halogenated cycloalkyl, halogenated heterocycloalkyl, substituted  
 15 alkyl, substituted cycloalkyl, substituted heterocycloalkyl, -NR<sub>11</sub>R<sub>11</sub>, -C(O)R<sub>11</sub>, -NO<sub>2</sub>, -C(O)NR<sub>11</sub>R<sub>11</sub>, -CN, -NR<sub>11</sub>C(O)R<sub>11</sub>, -S(O)<sub>2</sub>NR<sub>11</sub>R<sub>11</sub>, or -NR<sub>11</sub>S(O)<sub>2</sub>R<sub>11</sub>;

R<sub>13</sub> is selected from -OR<sub>11</sub>, -SR<sub>11</sub>, -NR<sub>11</sub>R<sub>11</sub>, -C(O)R<sub>11</sub>, -C(O)NR<sub>11</sub>R<sub>11</sub>, -CN, -NR<sub>11</sub>C(O)R<sub>11</sub>, -S(O)<sub>2</sub>NR<sub>11</sub>R<sub>11</sub>, -NR<sub>11</sub>S(O)<sub>2</sub>R<sub>11</sub>, -CF<sub>3</sub>, or -NO<sub>2</sub>;

R<sub>14</sub> is independently selected from -H, alkyl, halogenated alkyl, limited  
 20 substituted alkyl, cycloalkyl, halogenated cycloalkyl, substituted cycloalkyl, heterocycloalkyl, halogenated heterocycloalkyl, substituted heterocycloalkyl;

Each R<sub>18</sub> is independently selected from -H, alkyl, cycloalkyl, heterocycloalkyl, halogenated alkyl, halogenated cycloalkyl, halogenated heterocycloalkyl, substituted alkyl, substituted cycloalkyl, substituted  
 25 heterocycloalkyl, -OR<sub>11</sub>, -SR<sub>11</sub>, -NR<sub>11</sub>R<sub>11</sub>, -C(O)R<sub>11</sub>, -NO<sub>2</sub>, -C(O)NR<sub>11</sub>R<sub>11</sub>, -CN, -NR<sub>11</sub>C(O)R<sub>11</sub>, -S(O)<sub>2</sub>NR<sub>11</sub>R<sub>11</sub>, or -NR<sub>11</sub>S(O)<sub>2</sub>R<sub>11</sub>, -F, -Cl, -Br, or -I, or a bond directly or indirectly attached to the core molecule, provided that there is only one said bond to the core molecule within the 9-membered fused-ring moiety, further provided that the fused-ring moiety has 0-1 substituent selected from alkyl, cycloalkyl,  
 30 heterocycloalkyl, halogenated alkyl, halogenated cycloalkyl, halogenated heterocycloalkyl, substituted alkyl, substituted cycloalkyl, substituted heterocycloalkyl, -OR<sub>11</sub>, -SR<sub>11</sub>, -NR<sub>11</sub>R<sub>11</sub>, -C(O)R<sub>11</sub>, -NO<sub>2</sub>, -C(O)NR<sub>11</sub>R<sub>11</sub>, -CN, -NR<sub>11</sub>C(O)R<sub>11</sub>, -S(O)<sub>2</sub>NR<sub>11</sub>R<sub>11</sub>, or -NR<sub>11</sub>S(O)<sub>2</sub>R<sub>11</sub>, and further provided that the fused-

ring moiety has 0-3 substituent(s) selected from -F, -Cl, -Br, or -I; and pharmaceutically acceptable salts are useful to treat any one of or combination of schizophrenia, or psychosis.

Abbreviations which are well known to one of ordinary skill in the art may be used (e.g., "Ph" for phenyl, "Me" for methyl, "Et" for ethyl, "h" for hour or hours, min  
5 for minute or minutes, and "rt" for room temperature).

All temperatures are in degrees Centigrade.

Room temperature is within the range of 15-25 degrees Celsius.

Eq refers to equivalents.

10 Satd refers to saturated.

AChR refers to acetylcholine receptor.

nAChR refers to nicotinic acetylcholine receptor.

5HT<sub>3</sub>R refers to the serotonin-type 3 receptor.

FLIPR refers to a device marketed by Molecular Devices, Inc. designed to  
15 precisely measure cellular fluorescence in a high throughput whole-cell assay.  
(Schroeder et. al., *J. Biomolecular Screening*, 1(2), p 75-80, 1996).

TLC refers to thin-layer chromatography.

HPLC refers to high pressure liquid chromatography.

MeOH refers to methanol.

20 EtOH refers to ethanol.

IPA refers to isopropyl alcohol.

THF refers to tetrahydrofuran.

DMSO refers to dimethylsulfoxide.

DMF refers to dimethylformamide.

25 EtOAc refers to ethyl acetate.

TMS refers to tetramethylsilane.

TEA refers to triethylamine.

HATU refers to O-(7-azabenzotriazol-1-yl)-N,N,N', N'-tetramethyluronium  
hexafluorophosphate.

30 DIEA refers to N,N-diisopropylethylamine.

MLA refers to methyllycaconitine.

Ether refers to diethyl ether.

KH<sub>2</sub>PO<sub>4</sub> refers to potassium phosphate, monobasic.

NaClO<sub>2</sub> refers to sodium chlorite.

t-BuOH refers to tert-butanol.

Na<sub>2</sub>SO<sub>4</sub> refers to sodium sulfate.

MgSO<sub>4</sub> refers to magnesium sulfate.

5 K<sub>2</sub>CO<sub>3</sub> refers to potassium carbonate.

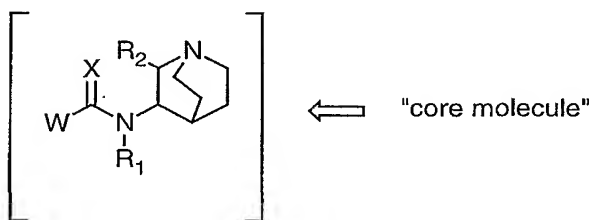
NH<sub>4</sub>OH refers to ammonium hydroxide.

NaHCO<sub>3</sub> refers to sodium bicarbonate.

CH<sub>3</sub>CN refers to acetonitrile.

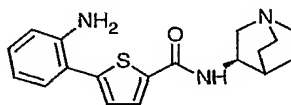
The ω carbon is determined by counting the longest carbon chain of the alkyl moiety with the C-1 carbon being the carbon attached to the W moiety of the core molecule and the ω carbon being the carbon furthest, e.g., separated by the greatest number of carbon atoms in the chain, from said C-1 carbon;

The core molecule is the quinuclidinyl-(carboxamide-type moiety)-W:



15 Therefore, when speaking of the ω carbon, the ω carbon is the carbon furthest from the core molecule and the C-1 carbon is the carbon attached to the core molecule by attachment to the W moiety of the core molecule.

One of the most conventionally accepted ways of naming the compound pictured below is 5-(2-aminophenyl)-N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]thiophene-2-carboxamide, but for one ordinarily skilled in the art, the following name also describes the same compound, N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenyl)-thiophene-2-carboxamide:



The two are used interchangeably in this patent.

25 The carbon atom content of various hydrocarbon-containing moieties is indicated by a prefix designating the minimum and maximum number of carbon atoms in the moiety, i.e., the prefix C<sub>i,j</sub> indicates a moiety of the integer 'i' to the

integer “j” carbon atoms, inclusive. Thus, for example, C<sub>1-6</sub> alkyl refers to alkyl of one to six carbon atoms, inclusive.

Halogen is F, Cl, Br, or I.

Alkyl is both straight and branched-chain moieties having from 1-6 carbon  
5 atoms.

Halogenated alkyl is an alkyl moiety having from 1-6 carbon atoms and having 1 to (2n+1) halogen atom(s) independently selected from -F, -Cl, -Br, or -I where n is the maximum number of carbon atoms in the moiety.

Substituted alkyl is an alkyl moiety having from 1-6 carbon atoms and having  
10 0-3 substituents independently selected from -F, -Cl, -Br, or -I, and further having 1 substituent selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -NO<sub>2</sub>, -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl.

Limited substituted alkyl is a substituted alkyl having from 1-6 carbon atoms  
15 and having 0-3 substituents independently selected from -F, -Cl, -Br, or -I, and further having 1 substituent on either only the ω carbon and selected from -OR<sub>11</sub>, -SR<sub>11</sub>, -NR<sub>11</sub>R<sub>11</sub>, -C(O)R<sub>11</sub>, -NO<sub>2</sub>, -C(O)NR<sub>11</sub>R<sub>11</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>11</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, or -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, or on any carbon with sufficient valency but not on the ω carbon and selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -NO<sub>2</sub>, -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN,  
20 -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl.

Alkenyl is straight and branched-chain moieties having from 2-6 carbon atoms and having at least one carbon-carbon double bond.

Halogenated alkenyl is an unsaturated alkenyl moiety having from 2-6 carbon atoms and having 1 to (2n-1) halogen atom(s) independently selected from -F, -Cl,  
25 -Br, or -I where n is the maximum number of carbon atoms in the moiety.

Substituted alkenyl is an unsaturated alkenyl moiety having from 2-6 carbon atoms and having 0-3 substituents independently selected from -F, or -Cl, and further having 1 substituent selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or  
30 substituted phenyl.

Limited substituted alkenyl is a substituted alkenyl having from 1-6 carbon atoms and having 0-3 substituents independently selected from -F, -Cl, -Br, or -I, and further having 1 substituent on either only the ω carbon and selected from -OR<sub>11</sub>,

-SR<sub>11</sub>, -NR<sub>11</sub>R<sub>11</sub>, -C(O)R<sub>11</sub>, -NO<sub>2</sub>, -C(O)NR<sub>11</sub>R<sub>11</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>11</sub>,  
 -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, or -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, or on any carbon with sufficient valency but not on  
 the ω carbon and selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -NO<sub>2</sub>,  
 -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or  
 5 substituted phenyl.

Alkynyl is straight and branched-chain moieties having from 2-6 carbon atoms  
 and having at least one carbon-carbon triple bond.

Halogenated alkynyl is an unsaturated alkynyl moiety having from 3-6 carbon  
 atoms and having 1 to (2n-3) halogen atom(s) independently selected from -F, -Cl,  
 10 -Br, or -I where n is the maximum number of carbon atoms in the moiety.

Substituted alkynyl is an unsaturated alkynyl moiety having from 3-6 carbon  
 atoms and having 0-3 substituents independently selected from -F, or -Cl, and further  
 having 1 substituent selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -CN,  
 -C(O)NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted  
 15 phenyl.

Limited substituted alkynyl is a substituted alkynyl having from 1-6 carbon  
 atoms and having 0-3 substituents independently selected from -F, -Cl, -Br, or -I, and  
 further having 1 substituent on either only the ω carbon and selected from -OR<sub>11</sub>,  
 -SR<sub>11</sub>, -NR<sub>11</sub>R<sub>11</sub>, -C(O)R<sub>11</sub>, -NO<sub>2</sub>, -C(O)NR<sub>11</sub>R<sub>11</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>11</sub>,  
 20 -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, or -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, or on any carbon with sufficient valency but not on  
 the ω carbon and selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -NO<sub>2</sub>,  
 -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or  
 substituted phenyl.

Cycloalkyl is a cyclic alkyl moiety having from 3-6 carbon atoms.

25 Halogenated cycloalkyl is a cyclic moiety having from 3-6 carbon atoms and  
 having 1-4 substituents independently selected from -F, or -Cl.

Substituted cycloalkyl is a cyclic moiety having from 3-6 carbon atoms and  
 having 0-3 substituents independently selected from -F, or -Cl, and further having 1  
 substituent selected from -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN,  
 30 -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, -NO<sub>2</sub>, phenyl, or substituted phenyl.

Heterocycloalkyl is a cyclic moiety having 4-7 atoms with 1-2 atoms within  
 the ring being -S-, -N(R<sub>3</sub>)- or -O-.

Halogenated heterocycloalkyl is a cyclic moiety having from 4-7 atoms with 1-2 atoms within the ring being -S-, -N(R<sub>3</sub>)-, or -O-, and having 1-4 substituents independently selected from -F, or -Cl.

Substituted heterocycloalkyl is a cyclic moiety having from 4-7 atoms with 1-2 atoms within the ring being -S-, -N(R<sub>3</sub>)-, or -O- and having 0-3 substituents independently selected from -F, or -Cl, and further having 1 substituent selected from -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -NO<sub>2</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl.

Substituted benzyl is a benzyl either having 1-4 substituents independently selected from -F, -Cl, -Br, or -I, or having 1 substituent selected from -R<sub>12</sub> and 0-3 substituents independently selected from -F, -Cl, -Br, or -I, provided that all substitution is on the phenyl ring of the benzyl.

Aryl is phenyl, substituted phenyl, naphthyl, or substituted naphthyl.

Substituted phenyl is a phenyl either having 1-4 substituents independently selected from -F, -Cl, -Br, or -I, or having 1 substituent selected from -R<sub>12</sub> and 0-3 substituents independently selected from -F, -Cl, -Br, or -I.

Substituted naphthyl is a naphthalene moiety either having 1-4 substituents independently selected from -F, -Cl, -Br, or -I, or having 1 substituent selected from -R<sub>12</sub> and 0-3 substituents independently selected from -F, -Cl, -Br, or -I, where the substitution can be independently on either the same ring or different rings of said naphthalene moiety.

Mammal denotes human and other mammals.

Compounds of the present invention may be in a form of pharmaceutically acceptable salts.

Brine refers to an aqueous saturated sodium chloride solution.

IR refers to infrared spectroscopy.

Lv refers to leaving groups within a molecule, including Br, Cl, OH, or mixed anhydride.

NMR refers to nuclear (proton) magnetic resonance spectroscopy, chemical shifts are reported in ppm (δ) downfield from TMS.

MS refers to mass spectrometry expressed as m/e or mass/charge unit. HRMS refers to high resolution mass spectrometry expressed as m/e or mass/charge unit.

M+H<sup>+</sup> refers to the positive ion of a parent plus a hydrogen atom. M-H<sup>-</sup> refers to the

negative ion of a parent minus a hydrogen atom.  $M+Na^+$  refers to the positive ion of a parent plus a sodium atom.  $M+K^+$  refers to the positive ion of a parent plus a potassium atom. EI refers to electron impact. ESI refers to electrospray ionization. CI refers to chemical ionization. FAB refers to fast atom bombardment.

5           Compounds of the present invention may be in the form of pharmaceutically acceptable salts. The term "pharmaceutically acceptable salts" refers to salts prepared from pharmaceutically acceptable non-toxic bases including inorganic bases and organic bases, and salts prepared from inorganic acids, and organic acids. Salts derived from inorganic bases include aluminum, ammonium, calcium, ferric, ferrous,  
10   lithium, magnesium, potassium, sodium, zinc, and the like. Salts derived from pharmaceutically acceptable organic non-toxic bases include salts of primary, secondary, and tertiary amines, substituted amines including naturally occurring substituted amines, cyclic amines, such as arginine, betaine, caffeine, choline, N, N-dibenzylethylenediamine, diethylamine, 2-diethylaminoethanol, 2-dimethylamino-  
15   ethanol, ethanolamine, ethylenediamine, N-ethylmorpholine, N-ethylpiperidine, glucamine, glucosamine, histidine, hydrabamine, isopropylamine, lysine, methylglucamine, morpholine, piperazine, piperidine, polyamine resins, procaine, purines, theobromine, triethylamine, trimethylamine, tripropylamine, and the like. Salts derived from inorganic acids include salts of hydrochloric acid, hydrobromic  
20   acid, hydroiodic acid, sulfuric acid, phosphoric acid, phosphorous acid and the like. Salts derived from pharmaceutically acceptable organic non-toxic acids include salts of  $C_{1-6}$  alkyl carboxylic acids, di-carboxylic acids, and tri-carboxylic acids such as acetic acid, propionic acid, fumaric acid, succinic acid, tartaric acid, maleic acid, adipic acid, and citric acid, and aryl and alkyl sulfonic acids such as toluene sulfonic  
25   acids and the like.

          By the term "effective amount" of a compound as provided herein is meant a nontoxic but sufficient amount of the compound(s) to provide the desired effect. As pointed out below, the exact amount required will vary from subject to subject, depending on the species, age, and general condition of the subject, the severity of the  
30   disease that is being treated, the particular compound(s) used, the mode of administration, and the like. Thus, it is not possible to specify an exact "effective amount." However, an appropriate effective amount may be determined by one of ordinary skill in the art using only routine experimentation.

The compounds of Formula I have optically active center(s) on the quinuclidine ring. Although it is desirable that the stereochemical purity be as high as possible, absolute purity is not required. This invention involves racemic mixtures and compositions of varying degrees of stereochemical purities. It is preferred to carry out stereoselective syntheses and/or to subject the reaction product to appropriate purification steps so as to produce substantially optically pure materials. Suitable stereoselective synthetic procedures for producing optically pure materials are well known in the art, as are procedures for purifying racemic mixtures into optically pure fractions.

The preferred compounds of the present invention have the R configuration at the C3 position of the quinuclidine ring. It is also preferred for the compounds of the present invention that X is O. Another group of compounds of Formula I includes compounds wherein X is O and R<sub>1</sub> is H. Another group of compounds of Formula I includes compounds wherein X is O and R<sub>2</sub> is H. Another group of compounds of Formula I includes compounds wherein X is O and R<sub>2</sub> is alkyl, halogenated alkyl, substituted alkyl, cycloalkyl, benzyl, substituted benzyl, or aryl.

The amount of therapeutically effective compound(s) that is administered and the dosage regimen for treating a disease condition with the compounds and/or compositions of this invention depends on a variety of factors, including the age, weight, sex and medical condition of the subject, the severity of the disease, the route and frequency of administration, and the particular compound(s) employed, and thus may vary widely. The compositions contain well known carriers and excipients in addition to a therapeutically effective amount of compounds of Formula I. The pharmaceutical compositions may contain active ingredient in the range of about 0.001-100 mg/kg/day for an adult, preferably in the range of about 0.1-50 mg/kg/day for an adult. A total daily dose of about 1-1000 mg of active ingredient may be appropriate for an adult. The daily dose can be administered in 1-4 doses per day.

In addition to the compound(s) of Formula I, the composition for therapeutic use may also comprise one or more non-toxic, pharmaceutically acceptable carrier materials or excipients. The term "carrier" material or "excipient" herein means any substance, not itself a therapeutic agent, used as a carrier and/or diluent and/or adjuvant, or vehicle for delivery of a therapeutic agent to a subject or added to a pharmaceutical composition to improve its handling or storage properties or to permit

or facilitate formation of a dose unit of the composition into a discrete article such as a capsule or tablet suitable for oral administration. Excipients can include, by way of illustration and not limitation, diluents, disintegrants, binding agents, adhesives, wetting agents, polymers, lubricants, glidants, substances added to mask or counteract a disagreeable taste or odor, flavors, dyes, fragrances, and substances added to improve appearance of the composition. Acceptable excipients include lactose, sucrose, starch powder, cellulose esters of alkanolic acids, cellulose alkyl esters, talc, stearic acid, magnesium stearate, magnesium oxide, sodium and calcium salts of phosphoric and sulfuric acids, gelatin, acacia gum, sodium alginate, polyvinylpyrrolidone, and/or polyvinyl alcohol, and then tableted or encapsulated for convenient administration. Such capsules or tablets may contain a controlled-release formulation as may be provided in a dispersion of active compound in hydroxypropylmethyl cellulose, or other methods known to those skilled in the art. For oral administration, the pharmaceutical composition may be in the form of, for example, a tablet, capsule, suspension or liquid. If desired, other active ingredients may be included in the composition.

In addition to the oral dosing, noted above, the compositions of the present invention may be administered by any suitable route, in the form of a pharmaceutical composition adapted to such a route, and in a dose effective for the treatment intended. The compositions may, for example, be administered parenterally, e.g., intravascularly, intraperitoneally, subcutaneously, or intramuscularly. For parenteral administration, saline solution, dextrose solution, or water may be used as a suitable carrier. Formulations for parenteral administration may be in the form of aqueous or non-aqueous isotonic sterile injection solutions or suspensions. These solutions and suspensions may be prepared from sterile powders or granules having one or more of the carriers or diluents mentioned for use in the formulations for oral administration. The compounds may be dissolved in water, polyethylene glycol, propylene glycol, ethanol, corn oil, cottonseed oil, peanut oil, sesame oil, benzyl alcohol, sodium chloride, and/or various buffers. Other adjuvants and modes of administration are well and widely known in the pharmaceutical art.

The serotonin type 3 receptor (5HT<sub>3</sub>R) is a member of a superfamily of ligand-gated ion channels, which includes the muscle and neuronal nAChR, the glycine receptor, and the  $\gamma$ -aminobutyric acid type A receptor. Like the other members of this

receptor superfamily, the 5HT<sub>3</sub>R exhibits a large degree of sequence homology with  $\alpha$  7 nAChR but functionally the two ligand-gated ion channels are very different. For example,  $\alpha$ 7 nAChR is rapidly inactivated, is highly permeable to calcium and is activated by acetylcholine and nicotine. On the other hand, 5HT<sub>3</sub>R is inactivated  
5 slowly, is relatively impermeable to calcium and is activated by serotonin. These experiments suggest that the  $\alpha$ 7 nAChR and 5HT<sub>3</sub>R proteins have some degree of homology, but function very differently. Indeed the pharmacology of the channels is very different. For example, Ondansetron, a highly selective 5HT<sub>3</sub>R antagonist, has little activity at the  $\alpha$ 7 nAChR. The converse is also true. For example, GTS-21, a  
10 highly selective  $\alpha$ 7 nAChR agonist, has little activity at the 5HT<sub>3</sub>R.

$\alpha$ 7 nAChR is a ligand-gated Ca<sup>++</sup> channel formed by a homopentamer of  $\alpha$ 7 subunits. Previous studies have established that  $\alpha$ -bungarotoxin ( $\alpha$ -btx) binds selectively to this homopentameric,  $\alpha$ 7 nAChR subtype, and that  $\alpha$ 7 nAChR has a high affinity binding site for both  $\alpha$ -btx and methyllycaconitine (MLA).  $\alpha$ 7 nAChR is  
15 expressed at high levels in the hippocampus, ventral tegmental area and ascending cholinergic projections from nucleus basalis to thalamocortical areas.  $\alpha$ 7 nAChR agonists increase neurotransmitter release, and increase cognition, arousal, attention, learning and memory.

Data from human and animal pharmacological studies establish that nicotinic  
20 cholinergic neuronal pathways control many important aspects of cognitive function including attention, learning and memory (Levin, E.D., *Psychopharmacology*, 108:417-31, 1992; Levin, E.D. and Simon B.B., *Psychopharmacology*, 138:217-30, 1998). For example, it is well known that nicotine increases cognition and attention in humans. ABT-418, a compound that activates  $\alpha$ 4 $\beta$ 2 and  $\alpha$ 7 nAChR, improves  
25 cognition and attention in clinical trials of Alzheimer's disease and attention-deficit disorders (Potter, A. et. al., *Psychopharmacology (Berl.)*, 142(4):334-42, Mar. 1999; Wilens, T. E. et. al., *Am. J. Psychiatry*, 156(12):1931-7, Dec. 1999). It is also clear that nicotine and selective but weak  $\alpha$ 7 nAChR agonists increase cognition and attention in rodents and non-human primates.

30 Schizophrenia is a complex multifactorial illness caused by genetic and non-genetic risk factors that produce a constellation of positive and negative symptoms. The positive symptoms include delusions and hallucinations and the negative symptoms include deficits in affect, attention, cognition and information processing.

No single biological element has emerged as a dominant pathogenic factor in this disease. Indeed, it is likely that schizophrenia is a syndrome that is produced by the combination of many low penetrance risk factors. Pharmacological studies established that dopamine receptor antagonists are efficacious in treating the overt  
5 psychotic features (positive symptoms) of schizophrenia such as hallucinations and delusions. Clozapine, an “atypical” antipsychotic drug, is novel because it is effective in treating both the positive and some of the negative symptoms of this disease. Clozapine’s utility as a drug is greatly limited because continued use leads to an increased risk of agranulocytosis and seizure. No other antipsychotic drug is effective  
10 in treating the negative symptoms of schizophrenia. This is significant because the restoration of cognitive functioning is the best predictor of a successful clinical and functional outcome of schizophrenic patients (Green, M.F., *Am J Psychiatry*, 153:321-30, 1996). By extension, it is clear that better drugs are needed to treat the cognitive disorders of schizophrenia in order to restore a better state of mental health to patients  
15 with this disorder.

One aspect of the cognitive deficit of schizophrenia can be measured by using the auditory event-related potential (P50) test of sensory gating. In this test, electroencephalographic (EEG) recordings of neuronal activity of the hippocampus are used to measure the subject’s response to a series of auditory “clicks” (Adler, L.E.  
20 et. al., *Biol. Psychiatry*, 46:8-18, 1999). Normal individuals respond to the first click with greater degree than to the second click. In general, schizophrenics and schizotypal patients respond to both clicks nearly the same (Cullum, C.M. et. al., *Schizophr. Res.*, 10:131-41, 1993). These data reflect a schizophrenic’s inability to “filter” or ignore unimportant information. The sensory gating deficit appears to be  
25 one of the key pathological features of this disease (Cadenhead, K.S. et. al., *Am. J. Psychiatry*, 157:55-9, 2000). Multiple studies show that nicotine normalizes the sensory deficit of schizophrenia (Adler, L.E. et. al., *Am. J. Psychiatry*, 150:1856-61, 1993). Pharmacological studies indicate that nicotine’s effect on sensory gating is via the  $\alpha 7$  nAChR (Adler, L.E. et. al., *Schizophr. Bull.*, 24:189-202, 1998). Indeed, the  
30 biochemical data indicate that schizophrenics have 50% fewer of  $\alpha 7$  nAChR receptors in the hippocampus, thus giving a rationale to partial loss of  $\alpha 7$  nAChR functionality (Freedman, R. et. al., *Biol. Psychiatry*, 38:22-33, 1995). Interestingly, genetic data indicate that a polymorphism in the promoter region of the  $\alpha 7$  nAChR gene is strongly

associated with the sensory gating deficit in schizophrenia (Freedman, R. et. al., *Proc. Nat'l Acad. Sci. USA*, 94(2):587-92, 1997; Myles-Worsley, M. et. al., *Am. J. Med. Genet*, 88(5):544-50, 1999). To date, no mutation in the coding region of the  $\alpha 7$  nAChR has been identified. Thus, schizophrenics express the same  $\alpha 7$  nAChR as  
5 non-schizophrenics.

Selective  $\alpha 7$  nAChR agonists may be found using a functional assay on FLIPR (see WO 00/73431 A2). FLIPR is designed to read the fluorescent signal from each well of a 96 or 384 well plate as fast as twice a second for up to 30 minutes. This assay may be used to accurately measure the functional pharmacology of  $\alpha 7$  nAChR  
10 and 5HT<sub>3</sub>R. To conduct such an assay, one uses cell lines that expressed functional forms of the  $\alpha 7$  nAChR using the  $\alpha 7/5$ -HT<sub>3</sub> channel as the drug target and cell lines that expressed functional 5HT<sub>3</sub>R. In both cases, the ligand-gated ion channel was expressed in SH-EP1 cells. Both ion channels can produce robust signal in the FLIPR assay.

15 The compounds of the present invention are  $\alpha 7$  nAChR agonists and may be used to treat a wide variety of diseases. For example, they may be used in treating schizophrenia, and psychosis.

Schizophrenia is a disease having multiple aspects. Currently available drugs are generally aimed at controlling the positive aspects of schizophrenia, such as  
20 delusions. One drug, Clozapine, is aimed at a broader spectrum of symptoms associated with schizophrenia. This drug has many side effects and is thus not suitable for many patients. Thus, there is a need for a drug to treat the cognitive and attention deficits associated with schizophrenia. Similarly, there is a need for a drug to treat the cognitive and attention deficits associated with schizoaffective disorders,  
25 or similar symptoms found in the relatives of schizophrenic patients.

Psychosis is a mental disorder characterized by gross impairment in the patient's perception of reality. The patient may suffer from delusions, and hallucinations, and may be incoherent in speech. His behavior may be agitated and is often incomprehensible to those around him. In the past, the term psychosis has been  
30 applied to many conditions that do not meet the stricter definition given above. For example, mood disorders were named as psychoses.

There are a variety of antipsychotic drugs. The conventional antipsychotic drugs include Chlorpromazine, Fluphenazine, Haloperidol, Loxapine, Mesoridazine,

Molindone, Perphenazine, Pimozide, Thioridazine, Thiothixene, and Trifluoperazine. These drugs all have an affinity for the dopamine 2 receptor.

These conventional antipsychotic drugs have several side effects, including sedation, weight gain, tremors, elevated prolactin levels, akathisia (motor  
5 restlessness), dystonia and muscle stiffness. These drugs may also cause tardive dyskinesia. Unfortunately, only about 70% of patients with schizophrenia respond to conventional antipsychotic drugs. For these patients, atypical antipsychotic drugs are available.

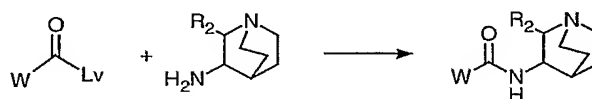
Atypical antipsychotic drugs generally are able to alleviate positive symptoms  
10 of psychosis while also improving negative symptoms of the psychosis to a greater degree than conventional antipsychotics. These drugs may improve neurocognitive deficits. Extrapyramidal (motor) side effects are not as likely to occur with the atypical antipsychotic drugs, and thus, these atypical antipsychotic drugs have a lower risk of producing tardive dyskinesia. Finally these atypical antipsychotic drugs cause  
15 little or no elevation of prolactin. Unfortunately, these drugs are not free of side effects. Although these drugs each produce different side effects, as a group the side effects include: agranulocytosis; increased risk of seizures, weight gain, somnolence, dizziness, tachycardia, decreased ejaculatory volume, and mild prolongation of QTc interval.

Finally, the compounds of the present invention may be used in combination  
20 therapy with typical and atypical anti-psychotic drugs. All compounds within the present invention are useful for and may also be used in combination with each other to prepare pharmaceutical compositions. Such combination therapy lowers the effective dose of the anti-psychotic drug and thereby reduces the side effects of the  
25 anti-psychotic drugs. Some typical anti-psychotic drugs that may be used in the practice of the invention include Haldol. Some atypical anti-psychotic drugs include Ziprasidone, Olanzapine, Risperidone, and Quetiapine.

Compounds of Formula I can be prepared as shown in Scheme 1. The key step  
30 in the preparation of this class of compounds is the coupling of commercially available 3-aminoquinuclidine ( $R_2 = H$ ) with the requisite acid chloride ( $L_v = Cl$ ), mixed anhydride (e.g.,  $L_v =$  diphenyl phosphoryl or acyloxy of the general formula of  $-O-C(O)-R_{L_v}$  where  $R_{L_v}$  includes phenyl or t-butyl), ester ( $L_v = OMe$  or  $OEt$ ), or

carboxylic acid ( $\text{Lv} = \text{OH}$ ) in the presence of an activating reagent. Suitable activating reagents are well known in the art, for examples see Kiso, Y.; Yajima, H. "Peptides" pp. 39-91, San Diego, CA, Academic Press, (1995), and include, but are not limited to, agents such as carbodiimides, phosphonium and uronium salts (such as uronium salt HATU).

Scheme 1



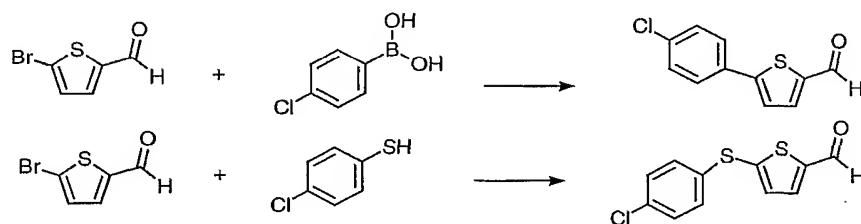
One of ordinary skill in the art will recognize that the methods described for the reaction of the unsubstituted 3-aminoquinuclidine ( $\text{R}_2 = \text{H}$ ) are equally applicable to substituted compounds ( $\text{R}_2 \neq \text{H}$ ). Such compounds can be prepared by reduction of the oxime of the corresponding 3-quinuclidinone (see *J. Labelled Compds. Radiopharm.*, 53-60 (1995) and *J. Med. Chem.* 988-995, (1998)). The oximes can be prepared by treatment of the 3-quinuclidinones with hydroxylamine hydrochloride in the presence of a base. The 3-quinuclidinones, where  $\text{R}_2 =$  substituted alkyl, cycloalkyl, substituted benzyl, can be prepared by known procedures (see *Tet. Lett.* 1015-1018, (1972), *J. Am. Chem. Soc.* 1278-1291 (1994), *J. Am. Chem. Soc.* 4548-4552 (1989), *Tetrahedron*, 1139-1146 (2000)). The 3-quinuclidinones, where  $\text{R}_2 =$  aryl, can be prepared by palladium catalyzed arylation as described in *J. Am. Chem. Soc.* 1473-1478 (1999) and *J. Am. Chem. Soc.* 1360-1370 (2000).

Preferably, when W is a thiophene or in some furan cases, the acid is converted into a mixed anhydride by treatment with diphenylchlorophosphate in the presence of TEA and  $\text{CH}_2\text{Cl}_2$  as the solvent. The resulting anhydride solution is directly reacted with aminoquinuclidine using aqueous DMF as the solvent. When W is furan, oxazole, oxadiazole, pyrrole, 5-thiazole, thiophene, or triazole, the acid is activated with a uronium salt, preferably HATU (see *J. Am. Chem. Soc.*, 4397 (1993)), in the presence of a base such as DIEA in DMF, and reacted directly with aminoquinuclidine to afford the desired amides. In the case where W is a 2-thiazole, 2-oxazole, or a thiadiazole, the amide bond is formed by the reaction of the amine and ester ( $\text{Lv} = \text{OEt}$ ) in an alcoholic solvent (see *Liebigs Ann. Chem.*, 1216-1231 (1980)).

It will be apparent to those skilled in the art that the requisite carboxylic acids can be obtained commercially or can be synthesized by known procedures. The

thiophene acids required in Examples 1-11, 13, and 41-42 can be synthesized from the corresponding aldehydes by oxidation with  $\text{NaClO}_2$  as described in *J. Chem. Soc. Perkin Trans. I.*, 789-794 (1999). The requisite aldehydes can be made as described in *J. Med. Chem.*, 1585-1599 (1997). An aryl boronic acid is reacted with a  
5 bromothiophene in the presence of a palladium (0) source, such as tetrakis-(triphenylphosphine)palladium (0), and a base, preferably aqueous sodium carbonate. The reaction works best if heated at reflux in THF/water for 24 hours. The thiophene acids of Examples 14-19 are prepared by similar methods as in Example 1 with modifications as described herein. The furan and thiophene acids required for  
10 Examples 20-30 are available commercially.

Scheme 2

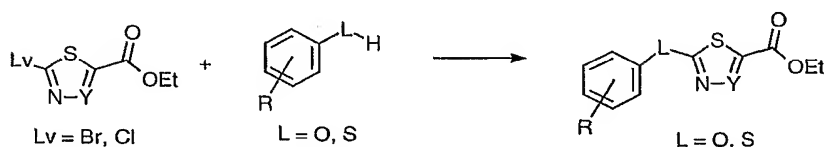


The thiophene acids for Examples 31-40 are synthesized from the corresponding esters by base catalyzed hydrolysis. Typical hydrolysis procedures are  
15 well known in the art. Preferably, the thiophene ester is treated with aqueous lithium hydroxide in a solvent such as dioxane. The esters are either commercially available or synthesized by reaction of a bromothiophene ester with the appropriate thiophenol or phenol as described in *Coll. Czech. Chem. Comm.*, 2360-2363 (1980). Namely, the sodium salt of the thiophenol or phenol is formed by treatment with a strong base like  
20 sodium hydride. The sodium salt is then reacted with a bromothiophene in a solvent such as acetone.

When W is thiazole, the required acids for Examples 44-49 are prepared by nucleophilic addition of the requisite phenol or thiophenol to 2-bromo-1,3-thiazole-5-carboxylic ethyl ester according to the procedure described in *Helv. Chim. Acta.*,  
25 2002-2022 (1997). Preferably, in EtOH utilizing  $\text{K}_2\text{CO}_3$  as a base (Scheme 3). The esters are hydrolyzed to the corresponding acids by procedures well known in the art. The 2-bromo-1,3-thiazole is prepared by the method described in *Roczniki Chemii Ann. Soc. Chim. Polonorum*, 1647-1658 (1972). The aryl 1,3-thiazole for Example 50 is prepared according to the procedure of Huntress and Pfister in *J. Am. Chem. Soc.*,

2167-2169 (1943). The requisite 1,3-thiazoles-5-acids for Examples 51 and 58-62 are commercially available. The 1,3-thiazole-5-acids required in Examples 52-57 can be synthesized from the corresponding esters by base hydrolysis via procedures well known in the art. The requisite esters can be prepared by a Suzuki reaction as  
 5 described in *J. Med. Chem.*, 4985-92 (1995). An aryl boronic acid is reacted with a bromothiazole ester in the presence of a palladium (0) source, such as tetrakis-(triphenylphosphine)palladium (0), and a base, preferably aqueous sodium carbonate. The 1,3,4-thiadiazole esters for Examples 64-72 are synthesized by nucleophilic  
 10 addition of the requisite phenol or thiophenol to 2-bromo-1,3,4-thiadiazole-5-carboxylic ethyl ester as described in *Can. J. Chem.*, 243-250 (1977) (Scheme 3).

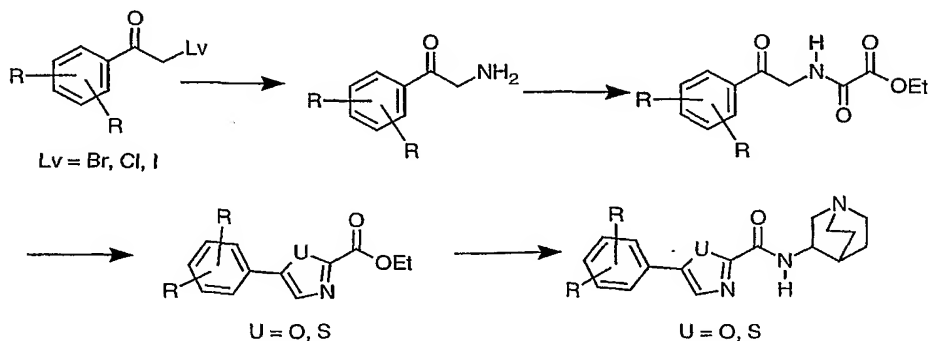
Scheme 3



The triazole for Example 75 and the oxadiazole for Example 76 are prepared by the methods of McKillop *et al.*, *Tetrahedron Lett.* 23, 3357-3360 (1982) with  
 15 modifications as described herein. The oxadiazole for Example 77 is prepared following the procedures of Snyder in *J. Org. Chem.* 3257-3269 (1990), Muchowski in *Can. J. Chem.* 3079-3082 (1972), and Crenshaw in US 4,001,238.

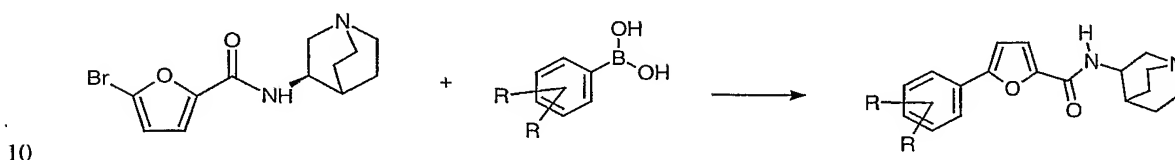
The 5-substituted-1,3-oxazole-2-esters for Examples 79-88 are synthesized according to procedures described in *J. Pharm. Soc. Japan*, 305-7 (1956) as shown in  
 20 Scheme 4. The 5-substituted-1,3-thiazole-2-esters for Examples 89-102 are synthesized according to procedures described in *Chem. Pharm. Bull.*, 4195-4198 (1982).

Scheme 4



The furans for Examples 103-130 are commercially available or can be prepared from their corresponding aldehydes or esters as described for the thiophenes (Examples 1-11). In the event that the furan is not commercially available, it can be prepared by the method of Bussolari and Rehborn described in *Org. Lett.* 965-7 (1999). Furan Examples 131-146 are prepared in a convergent means by a direct palladium catalyzed Suzuki coupling N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-furan-2-carboxamide with the requisite boronic acid by the method described in *Org. Lett.* 965-7 (1999), to yield directly the desired aryl amides (Scheme 5).

Scheme 5



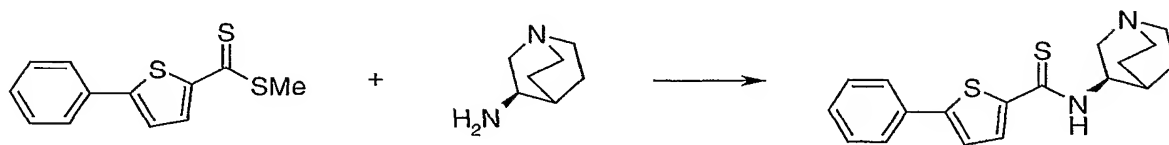
One of ordinary skill in the art would recognize that Examples 12, and 147-149 are prepared by reduction of the corresponding aryl nitro compounds by methods well known in the art, preferably by reduction with Pd/C in an alcoholic solvent such as EtOH under H<sub>2</sub>. The acid for example 150 is prepared by a Pd(0) catalyzed Sonogashira coupling of 5-bromo-2-fural and phenyl acetylene as described in *Tetrahedron Lett.*, 4467-70 (1975). The resulting aldehyde is converted to the desired analog by methods as described for Example 103. Example 151 is prepared by addition of the sodium salt of phenol to N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-furan-2-carboxamide. The requisite acid for Example 152 is prepared by bromination of methyl-5-bromo-1-methyl-1H-pyrrole-2-carboxylate, followed by similar Pd-catalyzed coupling as described for Example 1.

The 1,3-oxazole-2-carboxylic acid required for Example 153 is prepared by the method described in *J. Pharm. Sci. Japan* 305-7, (1956). 3-Phenyl-1,2,4-oxadiazole-5-carboxylic acid required for Example 154 is prepared by the method of Wurm as described in *Chem. Ber.*, 3133, (1889). The 2-phenyl-1,3-oxazole-5-carboxylic acid required for Example 155 is prepared by the method described in *Chem. Heterocycl. Compd.* (Engl. Transl.), 654-663, (1986). 2-Phenyl-1,3-oxazole-4-carboxylic acid required for Example 156 is prepared as described by Korte and Stoeriko, in *Chem. Ber.*, 1033-1042, (1960). The 5-phenylisoxazole-3-carboxylic acid

for Example 157 is prepared by the method of Vaughan and Spencer as described in *J. Org. Chem.* 1160-4, (1960).

Thioamides, such as Example 159, can be prepared from the requisite thioester by direct displacement of the thioester with aminoquinuclidine (Scheme 6). The thioester can be prepared as described in *J. Organometallic Chem.*, 95-98 (1987). One of ordinary skill in the art would quickly identify that compounds such as Example 159 could also be prepared directly from the amides exemplified throughout this patent by direct treatment with a reagent such as Lawesson's reagent (see Lawesson et. al. in *Bull. Soc. Chim. Belg.*, 229 (1978)) or  $P_4S_{10}$  (see *Chem. Rev.*, 45 (1961)).

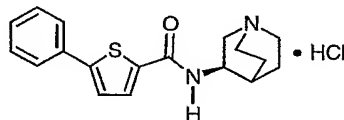
Scheme 6



The following examples are provided as examples and are not intended to limit the scope of this invention to only those provided examples and named compounds. Also, the salts made in the examples are only exemplary and are not intended to limit the invention. Any pharmaceutically acceptable salt can be made by one of ordinary skill in the art. Further, the naming of specific stereoisomers is for exemplification, and is not intended to limit in anyway the scope of the invention. The invention includes the following examples in pure stereoisomeric form or as racemic mixtures.

### Example 1

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenylthiophene-2-carboxamide hydrochloride:



Step 1a: Preparation of 5-phenylthiophene-2-carboxaldehyde.

5-Bromothiophene-2-carboxaldehyde (1.0g, 5.2 mmol) is added to a solution of tetrakis(triphenylphosphine)palladium(0) (180 mg, 0.16 mmol) in degassed THF

(10mL). The resulting solution is stirred for 5 minutes. A solution of phenylboronic acid (760 mg, 6.2 mmol) in THF (10mL) is added followed by aqueous Na<sub>2</sub>CO<sub>3</sub> (2M, 5.2mL). The mixture is heated at reflux for 24 hours. The reaction mixture is allowed to cool, poured into ether, and washed twice with water. The ether layer is dried over Na<sub>2</sub>SO<sub>4</sub> and concentrated *in vacuo*. The crude product is purified by flash column chromatography (1:1 hexanes:CH<sub>2</sub>Cl<sub>2</sub>) to yield the desired product (900 mg, 91%).  
<sup>1</sup>H NMR (300MHz, CDCl<sub>3</sub>) δ 7.38-7.45, 7.65-7.68, 7.73, 9.88.

Step 1b: Preparation of 5-phenylthiophene-2-carboxylic acid.

The product of Step 1a (750 mg, 4 mmol) is dissolved in a mixture of THF, t-BuOH, and water (2:1:1, 60 mL). KH<sub>2</sub>PO<sub>4</sub> (1.36 g, 10 mmol) is added followed by NaClO<sub>2</sub> (900 mg, 10 mmol). The mixture is stirred at room temperature for 5 days. Aqueous NaOH (2M, 10 mL) is added, and a majority of the organic solvents are removed *in vacuo* yielding an aqueous suspension. This suspension is diluted with water and washed three times with CH<sub>2</sub>Cl<sub>2</sub>. The aqueous layer is acidified to pH<6 with 25% H<sub>2</sub>SO<sub>4</sub> and the product is extracted three times with CH<sub>2</sub>Cl<sub>2</sub>. The combined organic washes are dried over Na<sub>2</sub>SO<sub>4</sub>, filtered and concentrated *in vacuo* to yield the desired product (417 mg, 51%). MS for C<sub>11</sub>H<sub>8</sub>O<sub>2</sub>S (ESI) (M-H)<sup>-</sup> *m/z* 203.

Step 1c Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenylthiophene-2-carboxamide hydrochloride.

TEA (210μL, 1.5mmol) is added to a suspension of the product of Step 1b (304mg, 1.5mmol) in CH<sub>2</sub>Cl<sub>2</sub> (5mL). Diphenylchlorophosphate (290μL, 1.4mmol) is added and the resulting solution is stirred at room temperature for 30 minutes. This solution is added to a solution of (R)-3-aminoquinuclidine dihydrochloride (279mg, 1.4mmol) and TEA (580μL, 4.2mmol) in DMF/water (5:1, 10mL). The resulting solution is allowed to stir overnight. MeOH is added and the mixture is poured through a column of AG50Wx2 ion exchange resin (H<sup>+</sup> form). The resin is washed with MeOH then the product is eluted with 5% TEA in MeOH. The eluent is evaporated to dryness. The hydrochloride salt is formed and crystallized from MeOH/IPA to yield the desired product (280mg, 57%). MS for C<sub>18</sub>H<sub>20</sub>N<sub>2</sub>OS (ESI) (M+H)<sup>+</sup> *m/z* 313.

The following compounds are made from the corresponding boronic acids according to the procedure of Example 1, making non-critical variations.

Example 2: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-thiophene-2-carboxamide hydrochloride (from 4-chlorophenylboronic acid). Yield 24% for 3  
5 steps. MS for  $C_{18}H_{19}ClN_2OS$  (ESI)  $(M+H)^+$   $m/z$  347.

Example 3: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-thiophene-2-carboxamide hydrochloride (from 3-chlorophenylboronic acid). Yield 16%. MS for  $C_{18}H_{19}ClN_2OS$  (ESI)  $(M+H)^+$   $m/z$  347.

Example 4: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-thiophene-2-  
10 carboxamide hydrochloride (from 2-chlorophenylboronic acid). Yield 48%. MS for  $C_{18}H_{19}ClN_2OS$  (ESI)  $(M+H)^+$   $m/z$  347.

Example 5: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2,3'-bithiophene-5-carboxamide hydrochloride (from 3-thiopheneboronic acid). Yield 49%. HRMS (FAB) calculated for  $C_{16}H_{18}N_2OS_2+H$  319.0939, found 319.0939.

15 Example 6: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-thiophene-2-carboxamide hydrochloride (from 2-nitrophenylboronic acid). Yield 45%. HRMS (FAB) calculated for  $C_{18}H_{19}N_3O_3S+H$  358.1225, found 358.1224.

Example 7: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-thiophene-2-carboxamide (from 3-nitrophenylboronic acid). Yield 25%. HRMS (FAB) calculated  
20 for  $C_{18}H_{19}N_3O_3S+H$  358.1225, found 358.1217.

Example 8: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-phenyl-thiophene-2-carboxamide hydrochloride (starting with phenyl boronic acid and 4-bromothiophene-2-carboxaldehyde in place of 5-bromothiophene-2-carboxaldehyde). Yield 35%. MS for  $C_{18}H_{20}N_2OS$  (ESI)  $(M+H)^+$   $m/z$  313.

25 Example 9: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-benzyloxyphenyl)-thiophene-2-carboxamide hydrochloride (from 3-benzyloxyphenylboronic acid). Yield 35%. MS for  $C_{25}H_{26}N_2O_2S$  (ESI)  $(M+H)^+$   $m/z$  419.

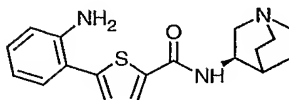
Example 10: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-benzyloxyphenyl)-thiophene-2-carboxamide hydrochloride (from 4-benzyloxyphenylboronic acid).  
30 Yield 34%. MS for  $C_{25}H_{26}N_2O_2S$  (ESI)  $(M+H)^+$   $m/z$  419.

Example 11: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoro-4-benzyloxyphenyl)-thiophene-2-carboxamide hydrochloride (from 3-fluoro-4-benzyloxyphenylboronic acid). Yield 41%. MS for  $C_{25}H_{25}FN_2O_2S$  (ESI)  $(M+H)^+$   $m/z$  436.

5

## Example 12

5-(2-Aminophenyl)-N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-thiophene-2-carboxamide:



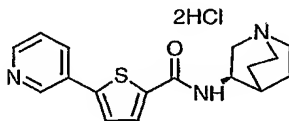
Step 12a: Preparation of 5-(2-aminophenyl)-N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]thiophene-2-carboxamide.

In a 200 mL Parr flask is placed the compound Example 6 (0.200 g, 0.51 mmol), EtOH (5 mL) and  $CH_2Cl_2$  (5 mL). The flask is shaken under 30 psi  $H_2$  for 16 hours. The contents are then filtered through a pad of celite with 10% MeOH- $CH_2Cl_2$  (150 mL). The solvents are removed *in vacuo*. This lot is then combined with a second lot prepared in a similar manner on the scale of 0.25 mmol. The product is re-dissolved in MeOH and loaded onto a column of AG50Wx2 ion exchange resin (H+ form). The resin is washed with MeOH then the product is eluted with 5% TEA in MeOH. The eluent is evaporated to dryness to yield the desired product (0.17 g, 69%). MS for  $C_{18}H_{21}N_3OS$  (EI)  $m/z$  (rel. intensity) 327 ( $M^+$ , 78), 327 (78), 203 (23), 202 (94), 130 (44), 125 (21), 117 (23), 109 (58), 96 (24), 83 (26), 70 (99).

20

## Example 13

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-pyridin-3-yl-thiophene-2-carboxamide dihydrochloride:



25

Step 13a: Preparation of 5-(3-pyridinyl)-2-thiophenecarboxaldehyde.

In a flask are placed 3-pyridinediethylborane (0.81 g, 5.5 mmol), 2-bromothiophene-5-carboxaldehyde (0.59 mL, 5.0 mmol), and Pd(PPh<sub>3</sub>)<sub>4</sub> (0.17g, 0.15 mmol). The flask is vacuum purged and nitrogen filled three times followed by addition of a 4:1 mixture of toluene-EtOH (8.3 mL) by syringe. After careful vacuum  
5 purge/nitrogen fill (3X), a solution of Na<sub>2</sub>CO<sub>3</sub> (2M, 5mL, 10.0 mmol) is added by syringe, and the flask is vacuum purged and nitrogen filled (3X). The reaction mixture is heated to 90°C and stirred for 22h. The reaction mixture is cooled to rt and diluted with H<sub>2</sub>O. The aqueous solution is extracted with ether (3X). The combined ether layers are then washed with water (3X) and brine (2X). The organic layer is  
10 dried over MgSO<sub>4</sub>, diluted with EtOAc for solubility, then filtered and concentrated. The crude product is then chromatographed over silica gel (10/30/50% EtOAc-heptane gradient) to provide the product as a yellow solid (0.34 g, 35%). NMR (400MHz, CDCl<sub>3</sub>): 9.97, 9.05, 8.68, 8.13, 7.83, 7.59-7.54.

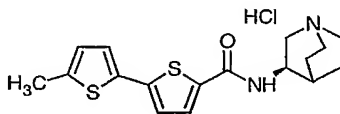
15 Step 13b: Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-pyridin-3-yl-thiophene-2-carboxamide dihydrochloride.

This compound is made from the product of Step 13a by using the procedure discussed in Steps 1b and 1c, making non-critical variations. The yield from 2 steps is 20%. MS for C<sub>17</sub>H<sub>19</sub>N<sub>3</sub>OS (ESI) (M+H)<sup>+</sup> *m/z* 314.

20

#### Example 14

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5'-methyl-2,2'-bithiophene-5-carboxamide hydrochloride:



25 Step 14a: Preparation of methyl 5'-methyl-2,2'-bithiophene-5-carboxylate.

In a flask are placed 2-bromothiophene-5-carboxylic acid (1.20 g, 5.81 mmol), 5-methylthiophene-2-boronic acid (0.99 g, 6.97 mmol), and Pd(PPh<sub>3</sub>)<sub>4</sub> (0.20 g, 0.17 mmol). The flask is then vacuum purged and nitrogen filled (3X). THF (12mL) is then added by syringe followed by vacuum purge and nitrogen fill (3X). A solution of

Na<sub>2</sub>CO<sub>3</sub> (2M, 5.8 mL, 11.6 mmol) is added followed by vacuum purge and nitrogen fill (3X). The reaction mixture is heated to reflux for 19h then cooled to rt and diluted with water. The aqueous solution is extracted with ether (3X). The aqueous layer is then acidified and extracted with EtOAc (3X). The combined organic layers are dried  
5 over MgSO<sub>4</sub>, filtered, and concentrated to provide an inseparable 3:1 mixture of 2-bromothiophene-5-carboxylic acid and bithiophene product. To separate the compounds the methyl esters are formed: In a flask are placed the aforementioned mixture, MeOH (50 mL) and conc H<sub>2</sub>SO<sub>4</sub> (5 drops) and heated to reflux for 24 h. The solution is concentrated and chromatographed over silica gel (2% acetone-heptane) to  
10 provide the product as a solid (0.37 g, 26% 2 steps). <sup>1</sup>H NMR (400MHz, CDCl<sub>3</sub>) δ 7.69, 7.10, 7.07, 6.73, 3.91, 2.52.

Step 14b: Preparation of 5'-methyl-2,2'-bithiophene-5-carboxylic acid.

The product of Step 14a (0.37 g, 1.54 mmol), dioxane (5 mL) and LiOH (1N,  
15 3.1 mL, 3.1 mmol) are placed in a flask. Additional dioxane (5 mL) is then added for solubility and stirred for 24 h at rt. 1N HCl is added slowly until pH<6, whereupon a precipitate forms. The precipitate is then collected by filtration, rinsed with water, and dried in a 70°C vacuum oven to provide the product as a yellow solid (0.30 g, 86%). MS for C<sub>10</sub>H<sub>8</sub>O<sub>2</sub>S<sub>2</sub> (ESI) (M-H)<sup>+</sup> *m/z* 223.

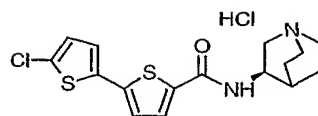
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Step 14c: Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5'-methyl-2,2'-bithiophene-5-carboxamide hydrochloride.

This compound is made using the product of Step 14b as the starting material and using the procedure discussed in Step 1c, making non-critical variations. Yield  
25 84%. MS C<sub>17</sub>H<sub>20</sub>N<sub>2</sub>OS<sub>2</sub> (EI) *m/z* (rel. intensity) 332 (M<sup>+</sup>, 90), 332 (90), 207 (99), 135 (90), 125 (55), 109 (86), 108 (43), 96 (52), 84 (52), 83 (48), 70 (88).

### Example 15

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5'-chloro-2,2'-bithiophene-5-  
30 carboxamide hydrochloride:



Step 15a: Preparation of 5'-chloro-2,2'-bithiophene-5-carboxaldehyde.

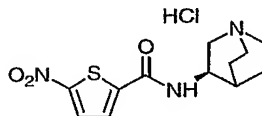
In a flask are placed 2-bromo-5-chlorothiophene (0.55 mL, 5.0 mmol) and Pd(PPh<sub>3</sub>)<sub>4</sub> (0.17 g, 0.15 mmol), and the flask is vacuum purged and nitrogen filled (3X). THF (10 mL) is added by syringe and stirred 10 min. In a separate flask is placed 5-formyl-2-thiopheneboronic acid (0.94 g, 6.0 mmol) and EtOH (2 mL) and stirred until dissolved. This mixture is added by syringe to the first flask followed by vacuum purge and nitrogen fill (3X). A solution of Na<sub>2</sub>CO<sub>3</sub> (2M, 5.0 mL, 10.0 mmol) is added by syringe followed by vacuum purge and nitrogen fill (3X). The reaction mixture is heated at 85°C for 20 h. The reaction is cooled to rt and diluted with water. The aqueous solution is extracted with ether (3X). The combined ether layers are washed with water (3X) then brine (2X). The ether is dried over MgSO<sub>4</sub>, filtered, and concentrated. The crude product is purified over silica gel (5% EtOAc-heptane) to provide the product as an orange solid (0.27 g, 24%). <sup>1</sup>H NMR (400MHz, CDCl<sub>3</sub>) δ 9.89, 7.69, 7.20, 7.16, 6.93.

Step 15b: Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5'-chloro-2,2'-bithiophene-5-carboxamide hydrochloride.

This compound is made from the product of Step 15a as the starting material by using the procedures discussed in Steps 1b and 1c, making non-critical variations. Yield for 2 steps 34%. HRMS (FAB) calculated for C<sub>16</sub>H<sub>17</sub>ClN<sub>2</sub>OS<sub>2</sub>+H 353.0549, found 353.0545.

#### Example 16

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-nitro-thiophene-2-carboxamide hydrochloride:

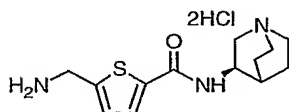


Step 16a: Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-nitrothiophene-2-carboxamide hydrochloride.

This compound is made from 2-nitrothiophene-5-carboxaldehyde by using the procedure discussed in Steps 1b and 1c, making non-critical variations. Yield for 2  
5 steps 48%. MS for  $C_{12}H_{15}N_3O_3S$  (EI)  $m/z$  (rel. intensity) 281 ( $M^+$ , 16), 156 (37), 125 (37), 110 (34), 109 (80), 96 (33), 84 (29), 82 (41), 81 (16), 70 (99), 55 (17).

### Example 17

5-(Aminomethyl)-N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-thiophene-2-  
10 carboxamide dihydrochloride:



Step 17a: Preparation of *tert*-butyl 2-thienylmethylcarbamate.

A 500 mL receiving flask is placed in an ice bath, and to the flask is added thiophene-2-methylamine (5.0 mL, 48.7 mmol),  $CH_2Cl_2$  (250 mL) then di-*tert*-butyl  
15 dicarbonate (12.7 g, 73.0 mmol) in 2-3 g portions over 5 min. The reaction mixture is stirred for 3 h then washed with 1N HCl (3X), 1N NaOH (3X) and brine (2X). The organic layer is dried over  $MgSO_4$ , filtered and concentrated to a yellow oil. The oil is chromatographed over silica gel (2.5-5% EtOAc-heptane gradient) to provide 8.70 g (84%) of the desired product as a clear oil. MS for  $C_{10}H_{15}NO_2S$  (ESI) ( $M+H$ ) $^+$   $m/z$   
20 214.

Step 17b: Preparation of 5-[(*tert*-butoxycarbonyl)amino]methyl}-2-thiophenecarboxylic acid.

In a flask is placed the product of Step 17a (3.50 g, 16.4 mmol) and dry THF  
25 (80 mL) then cooled in an acetone/solid  $CO_2$  bath. Lithium diisopropylamide (18.0 mL, 36.1 mmol, 2.0 M solution in heptane/THF/ethylbenzene) is added in a slow stream by syringe. The resulting orange solution is stirred for 10 min and then quenched with excess dry ice. The solution is warmed over 1h and the THF removed *in vacuo*. The crude product is diluted with  $CH_2Cl_2$  and washed with 1N HCl (3X).

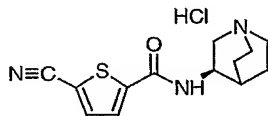
The  $\text{CH}_2\text{Cl}_2$  is removed *in vacuo* and replaced with EtOAc due to insolubility. The organic solution is dried over  $\text{MgSO}_4$ , filtered, and concentrated to provide a dark orange oil, which is chromatographed over silica gel (25-90% EtOAc-heptane gradient) to provide 0.873 g (21%) of the product as an off-white foam.  $^1\text{H}$  NMR (400MHz,  $\text{CDCl}_3$ )  $\delta$  12.96, 7.60, 7.56, 6.97, 4.28, 1.40.

Step 17c: Preparation of 5-(aminomethyl)-N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]thiophene-2-carboxamide dihydrochloride.

This compound is made by using the product of Step 17b as a starting material and using the procedure discussed in Step 1c, making non-critical variations. Yield 55%. MS for  $\text{C}_{13}\text{H}_{19}\text{N}_3\text{OS}$  (EI)  $m/z$  (rel. intensity) 265 ( $\text{M}^+$ , 6), 140 (24), 109 (79), 96 (24), 80 (31), 78 (99), 70 (80), 65 (28), 63 (99), 62 (39), 61 (77).

#### Example 18

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-cyano-thiophene-2-carboxamide hydrochloride:



Step 18a: Preparation of 5-cyano-thiophene-2-carboxylic acid.

In a 500 mL receiving flask are placed thiophene-2-carbonitrile (5.0 mL, 53.8 mmol) and THF (270 mL) and cooled in an acetone/solid  $\text{CO}_2$  bath. Lithium diisopropylamide (40.3 mL, 80.7 mmol, 2.0M solution in heptane/THF/ethylbenzene) is added in a slow stream via syringe. The solution is stirred for 10 min then quenched with an excess of dry ice. The reaction mixture is warmed in a water bath and the THF removed *in vacuo*. The slurry is taken up in 1N NaOH and extracted with ether (3X). The aqueous layer is then acidified to  $\text{pH} < 6$  with conc. HCl, whereupon a brown precipitate forms. This precipitate is filtered off and to the resulting eluent is added 1N HCl which results in precipitation of the product. The product is collected by filtration then triturated with  $\text{CH}_2\text{Cl}_2$ . Purification over silica gel (1:2.5:100 formic acid:MeOH: $\text{CH}_2\text{Cl}_2$ ) provides the product as a solid (1.79 g, 22%).  $^1\text{H}$  NMR (400MHz,  $\text{CDCl}_3$ )  $\delta$  14.10, 8.00, 7.80.

Step 18b: Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-cyano-thiophene-2-carboxamide hydrochloride.

This compound is made by using the product of Step 18a as a starting material and using the procedure discussed in Step 1c, making non-critical variations. Yield 57%. HRMS (FAB) calculated for  $C_{13}H_{15}N_3OS+H$  262.1014, found 262.1003.

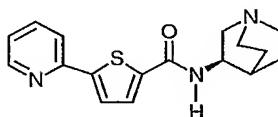
### Example 19

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-methoxy-thiophene-2-carboxamide hydrochloride:

This compound is prepared from 2-methoxythiophene according to the procedure used to make the compound of Example 18, making non-critical variations. Yield for 2 steps 33%. HRMS (FAB) calculated for  $C_{13}H_{18}N_2O_2S+H$  267.1167, found 267.1167.

### Example 20

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-2-yl)-thiophene-2-carboxamide:



Step 20a: TEA (50 $\mu$ L, 0.35mmol) is added to a suspension of 5-(pyridin-2-yl)-2-thiophenecarboxylic acid (72mg, 0.35mmol) in  $CH_2Cl_2$ :DMF (2:1, 1.5mL). Diphenylchlorophosphate (62 $\mu$ L, 0.3mmol) is added and the resulting solution is stirred at room temperature for 30 minutes. A solution of (R)-3-aminoquinuclidine (1M, 0.2mmol, 0.2mL) in DMF is added and the resulting solution is stirred overnight at room temperature. MeOH is added and the mixture is poured through a column of AG50Wx2 ion exchange resin ( $H^+$  form). The resin is washed with MeOH then the product is eluted with 5% TEA in MeOH. The eluent is evaporated to dryness to yield the desired product (15mg, 24%). MS for  $C_{17}H_{19}N_3OS$  (ESI) ( $M+H$ ) $^+$   $m/z$  314.

## Examples 21-30

The following compounds are made from the corresponding carboxylic acids according to the procedure of Example 20, making non-critical variations.

5 Example 21: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-[2,2']bithiophenyl-5-carboxamide (from 2,2'-bithiophene-5-carboxylic acid). Yield 13%. MS for  $C_{16}H_{18}N_2OS_2$  (ESI)  $(M+H)^+$   $m/z$  319.

Example 22: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(methylsulfanyl)-thiophene-2-carboxamide (from 5-methylsulfanylthiophene-2-carboxylic acid). Yield 84%. MS  
10 for  $C_{13}H_{18}N_2OS_2$  (ESI)  $(M+H)^+$   $m/z$  283.

Example 23: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-thiophene-2-carboxamide (from 5-chlorothiophene-2-carboxylic acid). Yield 6%. MS for  $C_{12}H_{15}ClN_2OS$  (ESI)  $(M+H)^+$   $m/z$  271.

Example 24: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-acetyl-thiophene-2-carboxamide (from 5-acetylthiophene-2-carboxylic acid). Yield 7%. MS for  
15  $C_{14}H_{18}N_2O_2S$  (ESI)  $(M+H)^+$   $m/z$  279.

Example 25: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-methyl-thiophene-2-carboxamide (from 5-methylthiophene-2-carboxylic acid). Yield 6%. MS for  $C_{13}H_{18}N_2OS$  (ESI)  $(M+H)^+$   $m/z$  251.

20 Example 26: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-thiophene-2-carboxamide (from 5-bromothiophene-2-carboxylic acid). Yield 8%. MS for  $C_{12}H_{15}BrN_2OS$  (ESI)  $(M+H)^+$   $m/z$  315/317.

Example 27: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylsulfanyl)-thiophene-2-carboxamide (from 5-phenylsulfanylthiophene-2-carboxylic acid). Yield 68%. MS  
25 for  $C_{18}H_{20}N_2OS_2$  (ESI)  $(M+H)^+$   $m/z$  345.

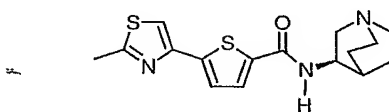
Example 28: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(acetyl-amino)-furan-2-carboxamide (from 5-acetylaminofuran-2-carboxylic acid). Yield 16%. MS for  $C_{14}H_{19}N_3O_3$  (ESI)  $(M+H)^+$   $m/z$  278.

Example 29: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-trifluoromethyl-furan-2-carboxamide (from 5-trifluoromethylfuran-2-carboxylic acid). Yield 11%. MS for  $C_{13}H_{15}F_3N_2O_2$  (ESI)  $(M+H)^+$   $m/z$  289.

Example 30: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methyl-5-trifluoromethyl-2H-pyrazole-3-yl)-thiophene-2-carboxamide (from 5-(2-methyl-5-trifluoromethyl-2H-pyrazole-3-yl)thiophene-2-carboxylic acid). Yield 9%. MS for  $C_{17}H_{19}F_3N_4OS$  (ESI)  $(M+H)^+$   $m/z$  385.

### Example 31

10 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylthiazol-4-yl)-thiophene-2-carboxamide:



Step 31a: Preparation of 5-(2-methylthiazol-4-yl)-thiophene-2-carboxylic acid.

Aqueous LiOH (1N, 1.5mL) is added to a solution of methyl 5-(2-methylthiazol-4-yl)-thiophene-2-carboxylate (81mg, 0.34mmol) in dioxane (1mL). The reaction is stirred at room temperature for 2 hours. Aqueous HCl (1N, 4mL) is added and the resultant precipitate is collected by filtration, washed with water, and dried to give the desired product (53mg, 69%).  $^1H$  NMR (300MHz, DMSO)  $\delta$  8.03, 7.69, 7.59, 2.70.

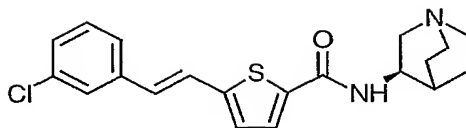
20 Step 31b: Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylthiazol-4-yl)-thiophene-2-carboxamide.

This compound is made by using the product of Step 31a as a starting material and using the procedure discussed in Step 20a, making non-critical variations. Yield 16%. MS for  $C_{16}H_{19}N_3OS_2$  (ESI)  $(M+H)^+$   $m/z$  334.

25

### Example 32

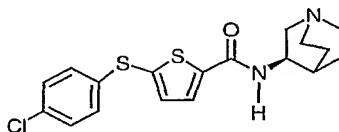
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(3-chlorophenyl)-vinyl]-thiophene-2-carboxamide:



This compound is prepared from methyl 5-[2-(3-chlorophenyl) vinyl]-thiophene-2-carboxylate according to the procedure used to make the compound of Example 31, making non-critical variations. Yield 19%. MS for  $C_{20}H_{21}ClN_2OS$  (ESI)  $(M+H)^+$   $m/z$  373.

### Example 33

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenylsulfanyl)-thiophene-2-carboxamide:



Step 33a: Preparation of 5-(4-chlorophenylsulfanyl)-thiophene-2-carboxaldehyde.

Sodium hydride (60%, 1.2g, 30mmol) is added to a solution of 4-chlorothiophenol (4.3g, 30mmol) in THF (30mL). The resulting solution is stirred for 10 minutes then the solvent is removed *in vacuo*. Acetone (60mL) is added followed by 5-bromothiophene-2-carboxaldehyde (3.0mL, 25mmol). The mixture is stirred at room temperature for 2 hours. The solvent is removed *in vacuo* and the resulting slurry diluted with  $CH_2Cl_2$ . This solution is washed three times with 1N NaOH then dried over  $MgSO_4$ , filtered and concentrated *in vacuo*. The crude product is purified by flash column chromatography (gradient of 1 to 5% EtOAc in heptane) to give the desired product (6.2g, 98%).  $^1H$  NMR (300MHz,  $CDCl_3$ )  $\delta$  7.13, 7.31-7.39, 7.63, 9.78.

Step 33b: Preparation of 5-(4-chlorophenylsulfanyl)-thiophene-2-carboxylic acid.

The product of Step 33a (6.1g, 24mmol) is dissolved in a mixture of THF, t-BuOH, and water (3:3:1, 255mL). 2-Methyl-2-butene (20.3mL, 192mmol) is added followed by  $KH_2PO_4$  (9.8g, 72mmol) and then  $NaClO_2$  (80%, 8.17g, 72.3mmol). The mixture is stirred at room temperature for 2 hours. Aqueous  $KHSO_4$  (0.5M, 200mL)

is added and the organic solvents are removed *in vacuo* to produce an aqueous suspension of the product. The precipitate is collected by filtration, dissolved in 1N NaOH and washed two times with ether. The aqueous solution is then acidified to pH<6 with concentrated HCl and a precipitate formed. The precipitate is collected by  
5 filtration and washed with 0.5M KHSO<sub>4</sub> then water. The solid is dried *in vacuo* to give the product (5.7g, 87%). MS for C<sub>11</sub>H<sub>7</sub>ClO<sub>2</sub>S<sub>2</sub> (ESI) (M-H)<sup>-</sup> *m/z* 269.

Step 33c: Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenylsulfanyl)-thiophene-2-carboxamide.

TEA (1.5mL, 10.9mmol) was added to a suspension of the product of Step 33b  
10 (2.7g, 10mmol) in CH<sub>2</sub>Cl<sub>2</sub> (100mL). Diphenylchlorophosphate (2.08mL, 10mmol) was added and the resulting solution is stirred at room temperature for 30 minutes. To this solution is added a solution of (R)-3-aminoquinuclidine (1M in DMF, 9.1mL, 9.1mmol). The resulting solution is stirred overnight at room temperature. MeOH is added and the mixture is poured through a column of AG50Wx2 ion exchange resin  
15 (H<sup>+</sup> form). The resin is washed with MeOH, then the product is eluted with 5% TEA in MeOH. The eluent is evaporated to dryness and the resulting solid is triturated with ether and dried *in vacuo* to yield the desired product (2.4g, 71%). Alternatively, these compounds can be converted to their hydrochloride salts and crystallized from MeOH/IPA. MS for C<sub>18</sub>H<sub>19</sub>ClN<sub>2</sub>OS<sub>2</sub> (ESI) (M+H)<sup>+</sup> *m/z* 379.

20

### Examples 34-39

The following compounds are made from the corresponding carboxylic acids according to the procedure of Example 33, making non-critical variations.

Example 34: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4 difluorophenyl-sulfanyl)-  
25 thiophene-2 carboxamide hydrochloride (from 2,4-difluorothiophenol). Yield for 3 steps 9%. MS for C<sub>18</sub>H<sub>18</sub>F<sub>2</sub>N<sub>2</sub>OS<sub>2</sub> (ESI) (M+H)<sup>+</sup> *m/z* 381.

Example 35: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl-sulfanyl)-thiophene-2-carboxamide (from 3-chlorothiophenol). Yield 50%. MS for C<sub>18</sub>H<sub>19</sub>ClN<sub>2</sub>OS<sub>2</sub> (ESI) (M+H)<sup>+</sup> *m/z* 379.

Example 36: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chloro-4-fluorophenylsulfanyl)-thiophene-2-carboxamide (from 3-chloro-4-fluorothiophenol). Yield 29%. MS for  $C_{18}H_{18}ClFN_2OS_2$  (ESI)  $(M+H)^+$   $m/z$  397.

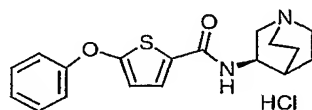
Example 37: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3-dichlorophenyl-sulfanyl)-thiophene-2-carboxamide hydrochloride (from 2,3-dichlorothiophenol). Yield 44%. MS for  $C_{18}H_{18}Cl_2N_2OS_2$  (ESI)  $(M+H)^+$   $m/z$  413.

Example 38: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4,5-trichlorophenyl-sulfanyl)-thiophene-2-carboxamide (from 2,4,5-trichlorothiophenol). Yield 53%. MS for  $C_{18}H_{17}Cl_3N_2OS_2$  (ESI)  $(M+H)^+$   $m/z$  449.

Example 39: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl-sulfanyl)-thiophene-2-carboxamide (from 3,4-dichlorothiophenol). Yield 21%. MS for  $C_{18}H_{18}Cl_2N_2OS_2$  (ESI)  $(M+H)^+$   $m/z$  413.

#### Example 40

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-thiophene-2-carboxamide hydrochloride:



Step 40a: 5-Phenoxy-thiophene-2-carboxaldehyde.

Phenol (3.3g, 35mmol) is added in portions to a suspension of 60% NaH (1.3g, 35mmol) in DMSO (100mL). The resulting mixture is stirred for 30 minutes then 5-nitrothiophene-2-carboxaldehyde (5g, 32mmol) is added. After 1 hour the reaction mixture is poured into water (1L) and washed with ether (4x500mL). The combined organic layers are dried over  $Na_2SO_4$ , filtered and evaporated to dryness. The resulting material is dissolved in MeOH and passed through a column (2.5cm x 20cm) of Amberjet 4400 ( $OH^-$  form). The eluent is dried *in vacuo* then evaporated twice from  $CH_3CN$ . The crude product is purified by column chromatography in EtOAc-hexanes (1:1) to yield the desired aldehyde (304mg, 5%).  $^1H$  NMR (300MHz,  $CDCl_3$ )  $\delta$  6.52, 7.20, 7.27, 7.45, 7.55, 9.75.

Step 40b: 5-Phenoxy-thiophene-2-carboxylic acid.

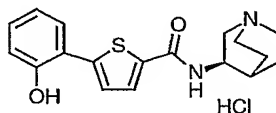
5-Phenoxy-thiophene-2-carboxaldehyde (325mg, 1.6mmol) is dissolved in a mixture of THF (10mL), t-BuOH (5mL) and water (5mL). NaH<sub>2</sub>PO<sub>4</sub> (650mg, 4.8mmol) is added followed by NaClO<sub>2</sub> (432mg, 4.8mmol). The resulting mixture is stirred for 24 hours at room temperature. Aqueous NaOH (2M, 5mL) is added, and the organic solvents are removed *in vacuo*. The resulting aqueous suspension is poured into water (50mL) and washed with ether (3x50mL). The aqueous layer is acidified to pH<2 with 25% H<sub>2</sub>SO<sub>4</sub> then washed with CH<sub>2</sub>Cl<sub>2</sub> (3x50mL). The combined organic washes are dried over Na<sub>2</sub>SO<sub>4</sub>, filtered and concentrated *in vacuo*. The crude product is dissolved in hot aqueous acetone and filtered. The solvents are gradually removed until a precipitate forms. The solid is collected by filtration and dried *in vacuo* to yield the desired product (192mg, 55%). MS for C<sub>11</sub>H<sub>7</sub>O<sub>3</sub>S (ESI) (M-H)<sup>+</sup> *m/z* 219.

Step 40c: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-thiophene-2-carboxamide hydrochloride.

The compound is made according to the procedure of Step 1c, starting with 5-phenoxy-thiophene-2-carboxylic acid and making non-critical variations. Yield (66%). MS for C<sub>18</sub>H<sub>21</sub>N<sub>2</sub>O<sub>2</sub>S (ESI) (M+H)<sup>+</sup> *m/z* 329.

#### Example 41

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-thiophene-2-carboxamide hydrochloride:



Step 41a: 5-(2-Hydroxyphenyl)-thiophene-2-carboxylic acid.

Tetrakis(triphenylphosphine)palladium(0) (133mg, 0.12mmol) is added to a solution of 5-bromothiophene-2-carboxylic acid (850mg, 4.1mmol) in degassed THF (10mL). The resulting solution is stirred for 5 minutes and then 2-(4,4,5,5-tetramethyl-1,3,2-dioxaborolan-2-yl) phenol (1g, 4.6mmol) is added followed by aqueous Na<sub>2</sub>CO<sub>3</sub> (2M, 6.9mL). The mixture is heated at reflux overnight. The reaction mixture is allowed to cool, poured into water (50mL), and washed with ether

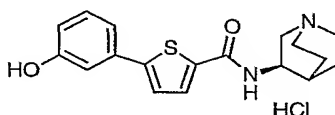
(3x50mL). The aqueous layer is acidified with concentrated HCl to pH<2. The resulting precipitate is collected by filtration, washed with water and dried *in vacuo* to yield the desired product (761mg, 83%). MS for C<sub>11</sub>H<sub>7</sub>O<sub>3</sub>S (ESI) (M-H)<sup>+</sup> *m/z* 219.

Step 41b: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-thiophene-2-carboxamide hydrochloride.

5-(2-Hydroxyphenyl)thiophene-2-carboxylic acid (650mg, 2.95mmol) is dissolved in DMF (30mL). (R)-3-Aminoquinuclidine dihydrochloride (587mg, 2.95mmol) is added followed by DIEA (1.6mL, 8.85mmol) and HATU (1.12g, 2.95mmol). The reaction is allowed to stir for 96 hours. The reaction mixture is diluted with MeOH and poured through a column of AG50Wx2. The column is washed with MeOH and then the product is eluted with 5% TEA in MeOH. The solvents are removed *in vacuo* and the resulting material is dissolved in CH<sub>3</sub>CN, concentrated and then dried *in vacuo*. The hydrochloride salt is formed and triturated with MeOH to yield the desired product (784mg, 73%). MS for C<sub>18</sub>H<sub>21</sub>N<sub>2</sub>O<sub>2</sub>S (ESI) (M+H)<sup>+</sup> *m/z* 329.

#### Example 42

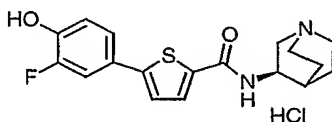
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-thiophene-2-carboxamide hydrochloride:



The compound is made from the 3-hydroxyphenylboronic acid according to the procedure of Example 41, making non-critical variations. Yield 46%. MS for C<sub>18</sub>H<sub>21</sub>N<sub>2</sub>O<sub>2</sub>S (ESI) (M+H)<sup>+</sup> *m/z* 329.

#### Example 43

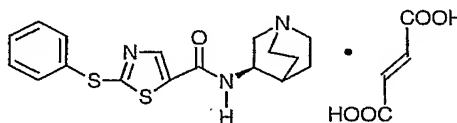
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoro-4-hydroxyphenyl)-thiophene-2-carboxamide hydrochloride:



Step 43a: The compound Example 11 (165mg, 0.38mmol) is dissolved in MeOH and poured through a plug of Amberjet 4400 (OH<sup>-</sup> form). The solvent is removed *in vacuo* and the product is redissolved in EtOH (2mL). This solution is added to a suspension of Pd/C (10%, 165mg) in EtOH (2mL). Cyclohexadiene (360mL, 3.8mmol) is added, and the reaction is heated at 60°C for 6 hours. The reaction mixture is diluted with MeOH and filtered through celite. The solvents are removed *in vacuo* then the hydrochloride salt is formed and crystallized from MeOH/CH<sub>3</sub>CN to yield the desired product (52mg, 36%). MS for C<sub>18</sub>H<sub>20</sub>FN<sub>2</sub>O<sub>2</sub>S (ESI) (M+H)<sup>+</sup> *m/z* 347.

#### Example 44

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(phenylsulfanyl)-1,3-thiazole-5-carboxamide fumarate:



Step 44a: Preparation of 2-phenylsulfanyl-thiazole-5-carboxylic acid ethyl ester.

A suspension of 2-bromo-1,3-thiazole-5-carboxylic acid ethyl ester (1.5 g, 6.15 mmol, 1 eq) and K<sub>2</sub>CO<sub>3</sub> (1.7 g, 12.3 mmol, 2 eq) in EtOH (60 mL) is cooled in an ice bath, and thiophenol (0.631 mL, 6.15 mmol, 1 eq) is added. The reaction is monitored by HPLC until the starting material is consumed. The reaction mixture is filtered (to remove a solid by-product), and the solvent is removed *in vacuo*. The crude mixture is purified by silica gel chromatography using a Biotage Flash 40S column using 2% EtOAc in hexanes to afford the product as an oil (0.784 g, 46%). MS (ESI) for C<sub>12</sub>H<sub>11</sub>NO<sub>2</sub>S<sub>2</sub> *m/z* 266.1 (M+H)<sup>+</sup>.

Step 44b: Preparation of 2-phenylsulfanyl-thiazole-5-carboxylic acid.

Potassium hydroxide (1.58 g, 28.2 mmol, 10 eq) is added to a solution of 2-phenylsulfanyl-thiazole-5-carboxylic acid ethyl ester (0.748 g, 2.82 mmol, 1 eq) in EtOH (15 mL) and water (10.5 mL). The reaction is stirred for 1.5 hours, diluted with water (30 mL) and EtOH (30 mL), and acidified by addition of 3 N HCl until a white precipitate forms. The precipitate is filtered and purified by recrystallization from

water and EtOH to give the product as a white crystalline solid (0.307 g). MS (ESI) for  $C_{10}H_7NO_2S_2$   $m/z$  235.9 (M-H)<sup>-</sup>.

Step 44c: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(phenylsulfanyl)-1,3-thiazole-5-carboxamide fumarate.

- 5 HATU (0.493 g, 1.30 mmol, 1 eq) is added to a solution of 2-phenylsulfanyl-thiazole-5-carboxylic acid (0.307 g, 1.30 mmol, 1 eq), (R)-(+)-3-aminoquinuclidine dihydrochloride (0.258 g, 1.30 mmol, 1 eq), and DIEA (0.677 mL, 3.89 mmol, 3 eq) in DMF (9.8 mL), the flask containing which was in an ice bath. The reaction is stirred overnight.  $CH_2Cl_2$  is added, and the mixture is washed with water three times.
- 10 The organic layer is dried over  $MgSO_4$ , and the solvent removed *in vacuo*. The product is purified using a Biotage Flash 40M column (1%  $NH_4OH$ /9%  $MeOH$ /90% $CH_2Cl_2$ ), affording a clear oil. The fumaric acid salt of the product is made, and recrystallized from MeOH and ether to give the product as a white crystalline solid (0.408 g, 72%). MS (ESI) for  $C_{17}H_{19}N_3OS_2$   $m/z$  346.1 (M+H)<sup>+</sup>.

15 Examples 45-49

The following compounds are prepared from the requisite phenols or thiophenols according to the procedures for Example 44, making non-critical variations.

- Example 45: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[(4-chlorophenyl)-sulfanyl]-  
20 1,3-thiazole-5-carboxamide fumarate (from 4-chlorothiophenol). Yield 39%. HRMS (FAB) calculated for  $C_{17}H_{18}ClN_3OS_2+H_1$  380.0658, found 380.0659.

Example 46: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-phenoxy-1,3-thiazole-5-carboxamide fumarate (from phenol). Yield 81%. HRMS (FAB) calculated for  $C_{17}H_{19}N_3O_2S+H_1$  330.1276, found 330.1269.

- 25 Example 47: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[(4-fluorophenyl)-sulfanyl]-1,3-thiazole-5-carboxamide fumarate (from 4-fluorothiophenol). Yield 43%. HRMS (FAB) calculated for  $C_{17}H_{18}FN_3OS_2+H_1$  363.0875, found 364.0945.

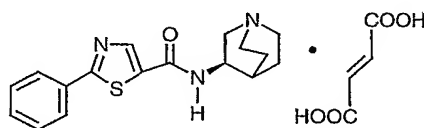
- Example 48: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(methylsulfanyl)-1,3-thiazole-5-carboxamide fumarate (from methanethiol). Yield 28%. HRMS (FAB) calculated  
30 for  $C_{12}H_{17}N_3OS_2+H_1$  284.0891, found 284.0894.

Example 49: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-chlorophenoxy)-1,3-thiazole-5-carboxamide fumarate (from 4-chlorophenol). Yield 42%. HRMS (FAB) calculated for  $C_{17}H_{18}ClN_3O_2S+H_1$  364.0886, found 364.0885.

5

## Example 50

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-phenyl-1,3-thiazole-5-carboxamide fumarate:



Step 50a: Preparation of 2-phenyl-thiazole-5-carboxylic acid ethyl ester.

10 A solution of  $\alpha$ -formyl- $\alpha$ -chloroacetate (9.34 g, 49.5 mmol, 1 eq) and thiobenzamide (6.79 g, 49.5 mmol, 1 eq) in EtOH (37.0 mL) is refluxed for 1 hour. The solution changes from an orange/brown color to a deep green. This solution is washed with water and extracted with  $CH_2Cl_2$ . The organic fraction is dried over  $Na_2SO_4$ , filtered, and the solvent removed *in vacuo*. The product is purified by  
15 column chromatography using a Biotage Flash 40M column (20% hexanes/EtOAc) to give the product as a deep orange oil (1.82 g, 15%). MS (ESI) for  $C_{12}H_{13}NO_3S$   $m/z$  252.1 (M+H)<sup>+</sup>.

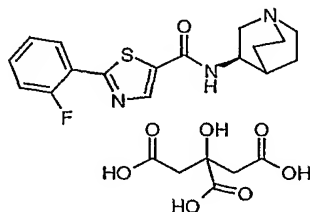
Step 50b: Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-phenyl-1,3-thiazole-5-carboxamide fumarate.

20 Making non-critical variations, the 2-phenyl-thiazole-5-carboxylic acid (1-azabicyclo[2.2.2]oct-3-yl)-amide is prepared using Steps 44b and 44c, starting with 2-phenyl-thiazole-5-carboxylic acid ethyl ester. The fumaric acid salt of the product is made and recrystallized from MeOH and ether to give the product as a white crystalline salt (77.4 mg, 22%). HRMS (FAB) calculated for  $C_{17}H_{19}N_3OS+H_1$   
25 314.1327, found 314.1336.

Example 51: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2,4-dimethyl-1,3-thiazole-5-carboxamide (from 2,3-dimethyl-thiazole-5-carboxylic acid as described in Step 1c). Yield 18%. HRMS (FAB) calculated for  $C_{13}H_{19}N_3OS+H_1$  266.1327, found 266.1332.

## Example 52

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-fluorophenyl)-1,3-thiazole-5-carboxamide citrate:



Step 52a: Preparation of ethyl 2-(2-fluorophenyl)-1,3-thiazole-5-carboxylate.

5 Tetrakis(triphenylphosphine)palladium (0) (0.58 g, 0.5 mmol), and a degassed solution of 2.0M Na<sub>2</sub>CO<sub>3</sub> (10 mL) are added to a degassed solution of 2-bromo-1,3-thiazole-5-carboxylic ethyl ester (1.18 g, 5.0 mmol) and 2-fluorophenylboronic acid (0.77 g, 5.5 mmol) in DME (10mL). The resulting suspension is stirred under argon at 80°C for 4 hours. The reaction mixture is cooled, diluted with EtOAc, and then  
10 washed with two portions of 1.0 M NaOH, then one portion of brine. The combined organic phases are concentrated *in vacuo*, and the resulting oil purified with flash chromatography. Yield 37%. HRMS (FAB) calculated for C<sub>12</sub>H<sub>10</sub>FNO<sub>2</sub>S+H 252.0495, found 252.0496.

Step 52b: Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-  
15 fluorophenyl)-1,3-thiazole-5-carboxamide citrate.

This compound is hydrolyzed and coupled according to Steps 44b and 44c, making non-critical variations. The citrate is prepared from the crude reaction mixture without chromatography, and crystallized. Yield 27%. HRMS (FAB) calculated for C<sub>17</sub>H<sub>18</sub>FN<sub>3</sub>OS+H 332.1233, found 332.1239.

20

### Examples 53-57

The following compounds are prepared from their requisite boronic acids according to the procedures for Example 52. The desired salt form is prepared directly from the crude reaction mixture without chromatography, and crystallized until the product is  
25 of analytical purity.

Example 53: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-fluorophenyl)-1,3-thiazole-5-carboxamide 4-methylbenzenesulfonate (from 3-fluorophenylboronic acid). Yield 20%. HRMS (FAB) calculated for  $C_{17}H_{18}FN_3OS+H$  332.1233, found 332.1225.

Example 54: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-fluorophenyl)-1,3-thiazole-5-carboxamide 4-methylbenzenesulfonate (from 4-fluorophenylboronic acid). Yield 51%. HRMS (FAB) calculated for  $C_{17}H_{18}FN_3OS+H$  332.1233, found 332.1239.

Example 55: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-hydroxyphenyl)-1,3-thiazole-5-carboxamide hydrochloride (from 2-(4,4,5,5-tetramethyl-1,3,2-dioxaborolan-2-yl)phenol). This compound was purified by reverse-phase preparative chromatography prior to the formation of the salt. Yield 0.5%. HRMS (FAB) calculated for  $C_{17}H_{19}N_3O_2S+H$  330.1276, found 330.1268.

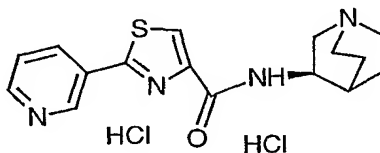
Example 56: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-methylphenyl)-1,3-thiazole-5-carboxamide hydrochloride (from 4-tolylphenylboronic acid). Yield 50%. MS (ESI) for  $C_{18}H_{21}N_3OS$   $m/z$  328.2 ( $M+H$ )<sup>+</sup>.

Example 57: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[4-(benzyloxy)phenyl]-1,3-thiazole-5-carboxamide hydrochloride (from (4-benzyloxyphenyl)boronic acid). Yield 98%. MS (ESI) for  $C_{24}H_{25}N_3O_2S$   $m/z$  420.3 ( $M+H$ )<sup>+</sup>.

#### Examples 58-62:

The following compounds are prepared from their requisite carboxylic acids according to the procedures from Step 44c. The desired salt form is prepared directly from the crude reaction mixture without chromatography, and crystallized.

Example 58: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-pyridin-3-yl-1,3-thiazole-4-carboxamide dihydrochloride (from (2-pyrid-3-yl)thiazole-4-carboxylic acid). HRMS (FAB) calculated for  $C_{16}H_{18}N_4OS+H$  315.1279, found 315.1289.



Example 59: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-2-phenyl-1,3-thiazole-5-carboxamide dihydrochloride (from 4-methyl-2-phenyl-1,3-thiazole-5-carboxylic acid). Yield 61%. HRMS (EI) calculated for  $C_{18}H_{21}N_3OS$  327.1405, found 327.1403.

Example 60: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-chlorophenyl)-4-methyl-1,3-thiazole-5-carboxamide 4-methylbenzenesulfonate (from 2-(4-chlorophenyl)-4-methyl-1,3-thiazole-5-carboxylic acid). Yield 85%. MS (ESI) for  $C_{18}H_{21}N_3OS$   $m/z$  328.2 (M+H)<sup>+</sup>.

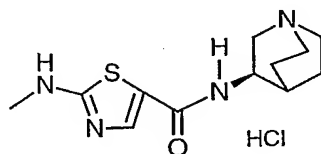
Example 61: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-2-pyridin-2-yl-1,3-thiazole-5-carboxamide hydrochloride (from 4-methyl-2-pyridin-2-yl-1,3-thiazole-5-carboxylic acid). Yield 57%. MS (ESI) for  $C_{18}H_{20}ClN_3OS$   $m/z$  362.2 (M+H)<sup>+</sup>.

Example 62: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-2-pyridin-4-yl-1,3-thiazole-5-carboxamide dihydrochloride (from 4-methyl-2-pyridin-4-yl-1,3-thiazole-5-carboxylic acid). Yield 87%. MS (ESI) for  $C_{17}H_{20}N_4OS$   $m/z$  329.2 (M+H)<sup>+</sup>.

15

## Example 63

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(methylamino)-1,3-thiazole-5-carboxamide hydrochloride:



Step 63a: Preparation of ethyl 2-[(tert-butoxycarbonyl)amino]-1,3-thiazole-5-carboxylate:

A flask is charged with a solution of ethyl 2-amino-1,3-thiazole-5-carboxylate (2.65 g, 15.4 mmol) and 4-dimethylaminopyridine (10 mg) in THF (75 mL). Di(tert-butyl) dicarbonate (3.6 mL, 15.4 mmol, 1.0 eq) and TEA (4.3 mL, 30.8 mmol, 2.0 eq) are added, and the resulting solution is stirred at room temperature for 90 minutes. The reaction mixture is concentrated to dryness, and the crude product is crystallized from  $CHCl_3$ /hexanes to give a light brown solid. Yield 68%. HRMS (FAB) calculated for  $C_{11}H_{16}N_2O_4S+H$  273.0909, found 273.0897.

Step 63b: Preparation of ethyl 2-[(tert-butoxycarbonyl)(methyl)amino]-1,3-thiazole-5-carboxylate.

A flask is charged with a suspension of sodium hydride (60% in mineral oil) (0.109 g, 2.72 mmol) in THF (5 mL). The ethyl 2-[(tert-butoxycarbonyl)amino]-1,3-thiazole-5-carboxylate (0.735 g, 2.70 mmol) is added, followed by iodomethane (175  $\mu$ L, 2.70 mmol) and the resulting suspension is heated to reflux for 3 hours, then cooled to room temperature. Water is added, followed by 1.0 N NaOH. The basic phase is extracted with 3 portions of EtOAc. The combined organic phases are washed with brine, dried over Na<sub>2</sub>SO<sub>4</sub>, filtered, and concentrated to give a clear oil purified with flash chromatography. Yield 45%. HRMS (FAB) calculated for C<sub>12</sub>H<sub>18</sub>N<sub>2</sub>O<sub>4</sub>S+H 287.1065, found 287.1068.

Step 63c: Preparation of 2-[(tert-butoxycarbonyl)(methyl)amino]-1,3-thiazole-5-carboxylic acid.

The product of Step 63b is hydrolyzed according to Step 44b, making non-critical variations. Yield 49%. HRMS (FAB) calculated for C<sub>10</sub>H<sub>14</sub>N<sub>2</sub>O<sub>4</sub>S+H 259.0752, found 259.0750.

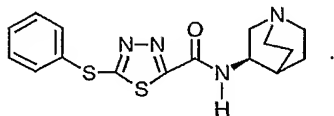
Step 63d: Preparation of tert-butyl 5-{[(3R)-1-azabicyclo[2.2.2]oct-3-ylamino]carbonyl}-1,3-thiazol-2-yl(methyl)carbamate.

This compound is coupled according to Step 44c. The citrate is prepared from the crude reaction mixture without chromatography, and crystallized until the product is of analytical purity. Yield 32%. MS (ESI) for C<sub>17</sub>H<sub>26</sub>N<sub>4</sub>O<sub>3</sub>S *m/z* 367.2 (M+H)<sup>+</sup>.

Step 63e: Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(methylamino)-1,3-thiazole-5-carboxamide hydrochloride.

A flask is charged with tert-butyl 5-{[(3R)-1-azabicyclo[2.2.2]oct-3-ylamino]carbonyl}-1,3-thiazol-2-yl(methyl)carbamate (72 mg, 0.18 mmol) in MeOH (5 mL). To this suspension is added a solution of 4.0N HCl/dioxane (10 mL) and the reaction mixture is stirred at room temperature for 3.5 hours. The solvents are removed *in vacuo* and the residue crystallized from IPA/ether. Yield 63%. MS (ESI) for C<sub>12</sub>H<sub>18</sub>N<sub>4</sub>OS *m/z* 267.2 (M+H)<sup>+</sup>.

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylsulfanyl)-1,3,4-thiadiazole-2-carboxamide:



Step 64a: Preparation of 5-phenylsulfanyl-[1,3,4]thiadiazole-2-carboxylic acid ethyl ester.

To a solution of 5-chloro-[1,3,4]thiadiazole-2-carboxylic acid ethyl ester (1.0 g, 5.2 mmol, 1 eq), and  $K_2CO_3$  (1.44 g, 10.4 mmol, 2eq) in  $CH_3CN$  (50 mL), thiophenol (0.53 mL, 5.2 mmol, 1 eq) is added. The reaction is stirred overnight at room temperature. The reaction is filtered to remove the salts, and the filtrate is washed with  $CH_3CN$ . The solvents are removed under reduced pressure, and the product is recrystallized from EtOH to give 5-phenylsulfanyl-[1,3,4]thiadiazole-2-carboxylic acid ethyl ester (1.03 g, 74%). MS (ESI) for  $C_{11}H_{10}N_2O_2S_2$   $m/z$  267.0 (M+H)<sup>+</sup>.

Step 64b: Preparation of 5-phenylsulfanyl-[1,3,4]thiadiazole-2-carboxylic acid.

The 5-phenylsulfanyl-[1,3,4]thiadiazole-2-carboxylic acid ethyl ester from Step 64a (1.0 g, 3.76 mmol, 1 eq) is dissolved in EtOH (40 mL), and cooled in an ice bath. To this, 2N NaOH (1.88 mL, 3.76 mmol, 1 eq) is added drop wise. The sodium salt of the acid precipitates out of solution. The reaction mixture is concentrated *in vacuo* to give a white crystalline product. The reaction mixture is carried through to the next step in its crude form to make the acid chloride. MS (ESI) for  $C_9H_6N_2S_2O_2$   $m/z$  237.0 (M-H)<sup>-</sup>.

Step 64c: Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylsulfanyl)-1,3,4-thiadiazole-2-carboxamide.

5-Phenylsulfanyl-[1,3,4]thiadiazole-2-carboxylic acid (3.76 mmol) is placed in a flask and put under nitrogen. The acid is chilled in an ice bath. Oxalyl chloride (9 mL) that has also been cooled in an ice bath is added drop wise to the acid. The excess oxalyl chloride is removed under reduced pressure. The acid chloride is dissolved in about 5 mL  $CH_2Cl_2$ . The freebase of (R)-3-aminoquinuclidine (0.5 g, 2.52 mmol, 0.67 eq) is also dissolved in  $CH_2Cl_2$  (5 mL), and the solution containing the acid chloride is cannulated into the flask containing the aminoquinuclidine. The

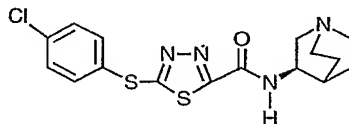
reaction is stirred overnight at room temperature. The reaction mixture is diluted with MeOH, and poured through a Dowex 50AGWX2 CH+ column. The impurities are removed by washing the column with MeOH, and the product is eluted by washing the column with a solution containing 5% TEA in MeOH. The solvent is removed under reduced pressure, and the product is further purified by silica gel chromatography using a Biotage Flash 40M column (5% MeOH/CH<sub>2</sub>Cl<sub>2</sub>). This gives Example 64 (0.289 g, 22%) as a tan oil. MS (ESI) for C<sub>16</sub>H<sub>18</sub>N<sub>4</sub>OS<sub>2</sub> *m/z* 347.2 (M+H)<sup>+</sup>.

- 10 Example 65: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-1,3,4-thiadiazole-2-carboxamide. This compound is prepared from phenol according to the procedures for Example 64, making non-critical variations. Yield 16%. HRMS (FAB) calculated for C<sub>16</sub>H<sub>17</sub>N<sub>3</sub>O<sub>3</sub>S 331.0991, found 331.1229 (M)<sup>+</sup>.

15

## Example 66

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-chlorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide:



- 5-(4-Chloro-phenylsulfanyl)-[1,3,4]thiadiazole-2-carboxylic acid ethyl ester (0.1 g, 0.33 mmol, 1 eq), the free base of (R)-(+)-3-aminoquinuclidine (0.041 g, 0.33 mmol, 1 eq), and EtOH (0.66 mL) are added to a vial, and the solution is heated at 80°C. The reaction is complete after 4 hours as determined by HPLC. The product is purified by silica gel chromatography using a Biotage Flash 40S column (0.5% NH<sub>4</sub>OH/9.5% MeOH/CH<sub>2</sub>Cl<sub>2</sub>) to give Example 66 (0.0254 g, 20%) as a pale oil.
- 25 HRMS (FAB) calculated for C<sub>16</sub>H<sub>17</sub>ClN<sub>4</sub>OS<sub>2</sub>+H<sub>1</sub> 381.0610, found 381.0617.

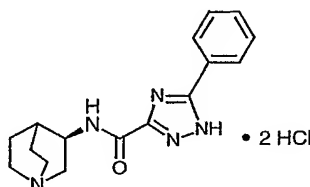
## Examples 67-72

The following compounds are prepared from the requisite phenols or thiophenols according to the procedures for Example 66, making non-critical variations.

- Example 67: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenoxy)-1,3,4-thiadiazole-2-carboxamide (from 4-fluorophenol). Yield 27%. HRMS (FAB) calculated for  $C_{16}H_{17}FN_4OS_2+H_1$  365.0906, found 365.0899.
- Example 68: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-1,3,4-thiadiazole-2-carboxamide (from 4-chlorophenol). Yield 18%. HRMS (FAB) calculated for  $C_{16}H_{17}ClN_4O_2S+H_1$  365.0839, found 365.0826.
- Example 69: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-fluorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide (from 3-fluorothiophenol). Yield 16%. HRMS (FAB) calculated for  $C_{16}H_{17}FN_4OS_2+H_1$  365.0906, found 365.0899.
- Example 70: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-chlorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide (from 2-chlorothiophenol). Yield 44%. HRMS (FAB) calculated for  $C_{16}H_{17}ClN_4OS_2+H_1$  381.0610, found 381.0625.
- Example 71: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-fluorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide (from 4-fluorothiophenol). Yield 34%. HRMS (FAB) calculated for  $C_{16}H_{17}FN_4OS_2+H_1$  365.0906, found 365.0921.
- Example 72: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-chlorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide (from 3-chlorothiophenol). Yield 34%. HRMS (FAB) calculated for  $C_{16}H_{17}ClN_4OS_2+H_1$  381.0610, found 381.0603.

### Example 75

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1H-1,2,4-triazole-3-carboxamide dihydrochloride:



Step 75a: Preparation of ethyl (2Z)-amino(benzoylhydrazono)ethanoate.

Ethyl thioamate (0.6g, 4.5mmol) and benzyl hydrazide 0.68g, 5.0mmol) in EtOH (20mL) are heated at reflux for 2 hours according to the procedure described in McKillop *et al.*, *Tetrahedron Lett.* 23, 3357-3360 (1982). The resulting solids are collected and washed with EtOH, and dried *in vacuo* to afford the desired product  
 5 (0.5g, 50%).

Step 75b: Preparation of ethyl 5-phenyl-1,3,4-oxadiazole-2-carboxylate and ethyl 5-phenyl-1H-1,2,4-triazole-3-carboxylate.

Modification of the procedure described in McKillop *et al.*, *Tetrahedron Lett.* 23, 3357-3360 (1982) affords a mixture of two products. In particular, heating ethyl  
 10 (2Z)-amino(benzoylhydrazono)ethanoate (0.5g, 2.1mmol) in mesitylene (15mL) at reflux overnight affords a clear solution that was concentrated *in vacuo*. The resultant material is purified by silica gel chromatography (20% EtOAc/hexanes) to afford a 2.2:1 mixture of ethyl 5-phenyl-1,3,4-oxadiazole-2-carboxylate (310mg, 67%) and ethyl 5-phenyl-1H-1,2,4-triazole-3-carboxylate (140mg, 30%), respectively. MS for  
 15  $C_{11}H_{10}N_2O_3$  (ethyl 5-phenyl-1,3,4-oxadiazole-2-carboxylate) (ESI)  $(M+H)^+$   $m/z$  219 and MS for  $C_{11}H_{10}N_2O_3$  (ethyl 5-phenyl-1H-1,2,4-triazole-3-carboxylate) (ESI)  $(M+H)^+$   $m/z$  218.

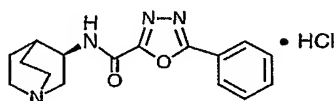
Step 75c: Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1H-1,2,4-triazole-3-carboxamide dihydrochloride.

20 This material is prepared using ethyl 5-phenyl-1H-1,2,4-triazole-3-carboxylate prepared in Step 75b and coupled to (R)-(+)-3-aminoquinuclidine. The product is purified as described in Example 66, making non-critical variations, to afford the desired product (13mg, 7.3%). MS for  $C_{16}H_{19}N_5O$  (ESI)  $(M+H)^+$   $m/z$  298.

25

### Example 76

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1,3,4-oxadiazole-2-carboxamide hydrochloride:

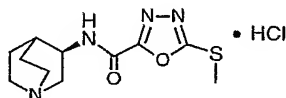


This material is prepared using ethyl 5-phenyl-1,3,4-oxadiazole-2-carboxylate  
 30 prepared in Step 75b and coupled to (R)-(+)-3-aminoquinuclidine and the product is

purified as described in Example 66, making non-critical variations. 7.3% Yield. MS for  $C_{16}H_{18}N_4O_2$  (ESI)  $(M+H)^+$   $m/z$  299.

### Example 77

5 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(methylthio)-1,3,4-oxadiazole-2-carboxamide hydrochloride:



Step 77a: Preparation of ethyl-1,3,4-oxadiazole-2-thione-5-carboxylate potassium salt dimethylsulfoxide solvate.

10 A mixture of ethyl hydrazino(oxo)acetate (3.7g, 28mmol), prepared as described by Benson, Gross, and Snyder in *J. Org. Chem.* 3257-3269 (1990), EtOH (25mL), carbon disulfide (8.5g, 6.6mL, 112mmol), DMSO (6mL) and a solution of KOH (1.57g, 28mmol) in water (2mL) are heated at reflux overnight, as described by Horning and Muchowski in *Can. J. Chem.* 3079-3082 (1972). The resultant solid is collected by vacuum filtration and washed to afford the desired product as a DMSO solvate that was used without further purification (8.2g of a semi-solid).

Step 77b: Preparation of methyl ethyl-1,3,4-oxadiazole-2-thione-5-carboxylate.

Following the procedure of Partyka and Crenshaw in US 4,001,238, a solution of ethyl-1,3,4-oxadiazole-2-thione-5-carboxylate potassium salt dimethylsulfoxide solvate (4.1g, 24mmol) and iodomethane (6.7g, 47mmol) are heated for 1 hour at reflux in EtOH (20mL). A solid is removed by vacuum filtration, and the filtrate is concentrated. Water (40mL) is added to the resultant material, a solid is isolated by vacuum filtration, dried at 50 °C/0.5 mmHg overnight to afford the desired product (2.3g, 50%).

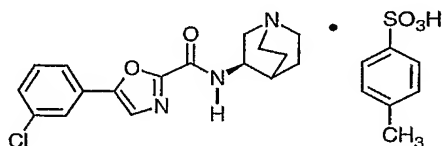
25 Step 77c: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(methylthio)-1,3,4-oxadiazole-2-carboxamide hydrochloride:

This material is prepared using methyl ethyl-1,3,4-oxadiazole-2-thione-5-carboxylate prepared in Step 77b and coupled to (R)-(+)-3-aminoquinuclidine as described in Example 66, making non-critical variations, to afford material that is

purified by silica gel chromatography (10% MeOH/CHCl<sub>3</sub> + 0.5% NH<sub>4</sub>OH). The hydrochloride salt is prepared upon crystallization of the desired product from 1M HCl/Et<sub>2</sub>O (65mg, 42%). MS for C<sub>11</sub>H<sub>16</sub>N<sub>4</sub>O<sub>2</sub>S (ESI) (M+H)<sup>+</sup> *m/z* 269.

### Example 79

- 5            *N*-[(3*R*)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1,3-oxazole-2-carboxamide 4-methylbenzenesulfonate:



Step 79a:        Preparation of 2-amino-1-(3-chlorophenyl)ethanone hydrochloride.

- A mixture of 3-chlorophenylacetyl bromide (5.18 g, 22.2 mmol, 1eq), sodium diformylimide (2.11 g, 22.2 mmol, 1 eq) and CH<sub>3</sub>CN (125 mL) is heated in an 80°C oil bath. After 3.5 h, the mixture is filtered and evaporated. The residue is treated with EtOH (40 mL) and HCl (10 mL, 12 N). The mixture was then heated in a 50°C water bath for 30 min and evaporated. The resulting solid is triturated with acetone and collected by filtration to afford the title compound (2.86 g, 62%). MS (ESI) for C<sub>8</sub>H<sub>8</sub>ClNO *m/z* 170 (M+H)<sup>+</sup>.
- 10
- 15

Step 79b:        Preparation of ethyl [[2-(3-chlorophenyl)-2-oxoethyl] amino](oxo) acetate.

- A mixture of 2-amino-1-(3-chlorophenyl)ethanone hydrochloride from Step 79a (2.83 g, 13.7 mmol, 1 eq), ethyl chlorooxoacetate (1.87 g, 13.7 mmol, 1 eq), and CH<sub>2</sub>Cl<sub>2</sub> (40 mL) is cooled in an ice-H<sub>2</sub>O bath. The mixture is treated with a solution of TEA (4.0 mL, 29 mmol, 2.1 eq) in CH<sub>2</sub>Cl<sub>2</sub> (20 mL), and the reaction is warmed to room temperature overnight. Water is added and the organic layer is separated, dried over MgSO<sub>4</sub>, filtered, and evaporated. The resulting solid is triturated with hexane/2-propanol and dried *in vacuo* to provide the title compound (2.70 g, 72%). MS (ESI) for C<sub>12</sub>H<sub>12</sub>ClNO<sub>4</sub> *m/z* 270 (M+H)<sup>+</sup>.
- 20
- 25

Step 79c:        Preparation of ethyl 5-(3-chlorophenyl)-1,3-oxazole-2-carboxylate.

A mixture of the product from Step 79b (1.28 g, 4.70 mmol, 1 eq), benzene (8 mL), and POCl<sub>3</sub> (2.0 mL, 21 mmol) is heated under reflux for 65 h and cooled. The

mixture is then evaporated and extracted between  $\text{CHCl}_3$  and water. The organic layer is separated, dried over  $\text{MgSO}_4$ , filtered, and evaporated. The residue is crystallized from EtOH to give the title product (0.61 g, 51%). MS (ESI) for  $\text{C}_{12}\text{H}_{10}\text{ClNO}_3$   $m/z$  252 ( $\text{M}+\text{H}$ )<sup>+</sup>.

- 5 Step 79d: Preparation of *N*-[(3*R*)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1,3-oxazole-2-carboxamide 4-methylbenzenesulfonate.

A mixture of the product from Step 79c (0.419 g, 1.66 mmol, 1.0 eq), (*R*)-(+)-3-aminoquinuclidine (0.221 g, 1.75 mmol, 1.05 eq) and EtOH (4.0 mL) is heated under reflux. After 64 h, the mixture is cooled and the contents are filtered through  
10 glass wool into *p*-toluenesulfonic acid monohydrate (0.315g, 1.65 mmol, 1.0 eq). The resulting solid is filtered and dried *in vacuo* to afford the title compound (0.59 g, 65%). HRMS (FAB) calculated for  $\text{C}_{17}\text{H}_{18}\text{ClN}_3\text{O}_2+\text{H}_1$  332.1165, found 332.1180.

#### Examples 80-88

- 15 The following compounds are prepared from the requisite acylhalides according to the procedures for Example 79, making non-critical variations. For some examples, the product of Step 79a is commercially available.

Example 80: *N*-[(3*R*)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-1,3-oxazole-2-carboxamide 4-methylbenzenesulfonate (from 2-bromo-1-(3-methoxyphenyl)ethanone). Yield 20%. HRMS (FAB) calculated for  $\text{C}_{18}\text{H}_{21}\text{N}_3\text{O}_3+\text{H}_1$   
20 328.1661, found 328.1660.

Example 81: *N*-[(3*R*)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-1,3-oxazole-2-carboxamide 4-methylbenzenesulfonate (from 2-amino-1-(3-nitrophenyl)ethanone hydrochloride). Yield 8%. HRMS (FAB) calculated for  $\text{C}_{17}\text{H}_{18}\text{N}_4\text{O}_4+\text{H}_1$  343.1406,  
25 found 343.1405.

Example 82: *N*-[(3*R*)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-1,3-oxazole-2-carboxamide 4-methylbenzenesulfonate (from 2-amino-1-(4-methoxyphenyl)ethanone hydrochloride). Yield 23%. HRMS (FAB) calculated for  $\text{C}_{18}\text{H}_{21}\text{N}_3\text{O}_3+\text{H}_1$  328.1661, found 328.1662.

Example 83: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-1,3-oxazole-2-carboxamide 4-methylbenzenesulfonate (from 2-bromo-1-(2-nitrophenyl)ethanone). Yield 4%. HRMS (FAB) calculated for  $C_{17}H_{18}N_4O_4+H_1$  343.1406, found 343.1405.

Example 84: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1,3-oxazole-2-carboxamide 4-methylbenzenesulfonate (from 2-bromo-1-(2-methoxyphenyl)ethanone). Yield 13%. HRMS (FAB) calculated for  $C_{18}H_{21}N_3O_3+H_1$  328.1661, found 328.1661.

Example 85: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1,3-oxazole-2-carboxamide 4-methylbenzenesulfonate (from 2-bromo-1-(4-fluorophenyl)ethanone). Yield 5%. HRMS (FAB) calculated for  $C_{17}H_{18}FN_3O_2+H_1$  316.1461, found 316.1470.

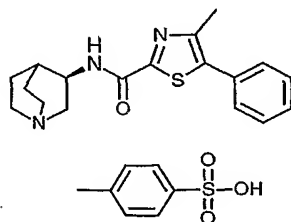
Example 86: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-1,3-oxazole-2-carboxamide 4-methylbenzenesulfonate (from 2-bromo-1-(2-chlorophenyl)ethanone). Yield 10%. HRMS (FAB) calculated for  $C_{17}H_{18}ClN_3O_2+H_1$  332.1165, found 332.1168.

Example 87: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-1,3-oxazole-2-carboxamide 4-methylbenzenesulfonate (from 3-(bromoacetyl)-benzonitrile). Yield 3%. HRMS (FAB) calculated for  $C_{18}H_{21}N_3O_3+H_1$  323.1508, found 323.1516.

Example 88: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-1,3-oxazole-2-carboxamide 4-methylbenzenesulfonate (from 2-amino-1-(4-bromophenyl)ethanone hydrochloride). Yield 6%. HRMS (FAB) calculated for  $C_{17}H_{18}BrN_3O_2+H_1$  376.0661, found 376.0660.

#### Example 89

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-5-phenyl-1,3-thiazole-2-carboxamide 4-methylbenzenesulfonate:



Step 89a: Preparation of 2-amino-1-phenylpropan-1-one hydrochloride.

2-Bromo-1-phenylpropan-1-one (8.97g, 42.1mmol, 1eq) is added dropwise to a suspension of diformylimide sodium salt (4.80g, 50.5mmol, 1.2eq) in 80mL CH<sub>3</sub>CN. The reaction is stirred for 60h at 70-75°C. The hot mixture is filtered to  
5 remove the salts and the solids are washed with CH<sub>3</sub>CN. The combined filtrates are concentrated *in vacuo*, dissolved in 40mL 6N HCl and heated under reflux for 0.75h. The solvents are removed under reduced pressure and the product is recrystallized from IPA to give 2-amino-1-phenylpropan-1-one hydrochloride (6.15g, 79%). MS (ESI) for C<sub>9</sub>H<sub>11</sub>NO *m/z* 150.2 (M+H)<sup>+</sup>.

10 Step 89b: Preparation of ethyl [(1-methyl-2-oxo-2-phenylethyl)amino] (oxo)acetate.

TEA (3.22mL, 0.0231mol, 2.1eq) is added dropwise to a suspension of the product from Step 89a (2.05g, 11.0mmol, 1eq) and ethyl oxalyl chloride (1.24mL, 11.0mmol, 1eq) in 50mL CH<sub>2</sub>Cl<sub>2</sub> in an ice/water bath. The mixture is allowed to  
15 slowly warm to room temperature. After stirring overnight, water and 20mL 1N HCl are added. The aqueous layer is extracted with CH<sub>2</sub>Cl<sub>2</sub>. The combined organic layers are dried over MgSO<sub>4</sub>, filtered and concentrated to give ethyl [(1-methyl-2-oxo-2-phenylethyl)amino](oxo)acetate as a yellow oil (2.58g, 94%). MS (ESI) for C<sub>13</sub>H<sub>15</sub>NO<sub>4</sub> *m/z* 250.2 (M+H)<sup>+</sup>.

20 Step 89c: Preparation of ethyl 4-methyl-5-phenyl-1,3-thiazole-2-carboxylate.

The product from Step 89b (2.58g, 10.4mmol, 1eq) and P<sub>2</sub>S<sub>5</sub> (4.83g, 10.9mmol, 1.05eq) are suspended in 30mL CHCl<sub>3</sub>. The mixture is heated under reflux. After 12h, water and solid K<sub>2</sub>CO<sub>3</sub> are carefully added until all material dissolves. The aqueous layer is made sufficiently basic with 1N NaOH (pH more than  
25 10) and extracted with EtOAc. The combined organic layers are washed with 1N NaOH and brine, dried over MgSO<sub>4</sub>, filtered and concentrated to give ethyl 4-methyl-5-phenyl-1,3-thiazole-2-carboxylate as a yellow oil (2.51g, 98%). MS (ESI) for C<sub>13</sub>H<sub>13</sub>NO<sub>2</sub>S *m/z* 248.1 (M+H)<sup>+</sup>.

Step 89d: Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-5-  
30 phenyl-1,3-thiazole-2-carboxamide 4-methylbenzenesulfonate.

The freebase of (*R*)-3-aminoquinuclidine (1.0g, 5.02mmol, 2.26eq) dissolved in THF (~5mL) is added to a solution of the product from Step 89c (0.55g, 2.22mmol, 1eq) in 10mL EtOH. The mixture is heated under reflux. After 48h, the mixture is cooled and concentrated *in vacuo*. The residue is purified by silica gel

5 chromatography using a Biotage Flash 40S column (90:9:1 CHCl<sub>3</sub>/MeOH/NH<sub>4</sub>OH) to provide N-[(3*R*)-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-5-phenyl-1,3-thiazole-2-carboxamide (0.40g, 55%) as a yellow oil. The *p*-toluenesulfonic acid salt of the product is made and recrystallized from IPA/ether to give the product as light yellow solid. MS (ESI) for C<sub>18</sub>H<sub>21</sub>N<sub>3</sub>OS m/z 328.2 (M+H)<sup>+</sup>.

10

## Examples 90-102

The following compounds are prepared from the requisite ethyl 5-substituted-1,3-thiazole-2-carboxylates according to the procedures outlined in Example 79 or Example 66 (thiadiazole amide), making non-critical variations.

Example 90: N-[(3*R*)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1,3-thiazole-2-  
15 carboxamide (from 2-amino-1-phenylethanone hydrochloride). Yield 36%. HRMS (FAB) calculated for C<sub>17</sub>H<sub>19</sub>N<sub>3</sub>OS+H<sub>1</sub> 314.1327, found 314.1330.

Example 91: N-[(3*R*)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-1,3-thiazole-2-carboxamide 4-methylbenzenesulfonate (from 2-amino-1-(4-bromophenyl)-ethanone hydrochloride). Yield 18%. HRMS (FAB) calculated for  
20 C<sub>17</sub>H<sub>18</sub>BrN<sub>3</sub>OS+H<sub>1</sub> 392.0432, found 392.0423.

Example 92: N-[(3*R*)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-1,3-thiazole-2-carboxamide 4-methylbenzenesulfonate (from 2-amino-1-(3-nitrophenyl)-ethanone hydrochloride). Yield 47%. HRMS (FAB) calculated for C<sub>17</sub>H<sub>18</sub>N<sub>4</sub>O<sub>3</sub>S+H<sub>1</sub>  
359.1178, found 359.1165.

25 Example 93: N-[(3*R*)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-1,3-thiazole-2-carboxamide 4-methylbenzenesulfonate (from 2-bromo-1-(3-methoxyphenyl)-ethanone). Yield 14%. HRMS (FAB) calculated for C<sub>18</sub>H<sub>21</sub>N<sub>3</sub>O<sub>2</sub>S+H<sub>1</sub> 344.1432, found 344.1423.

Example 94: N-[(3*R*)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-1,3-thiazole-  
30 2-carboxamide 4-methylbenzenesulfonate (from 2-bromo-1-(2-chlorophenyl)-

ethanone). Yield 28%. HRMS (FAB) calculated for  $C_{17}H_{18}ClN_3OS+H_1$  348.0937, found 348.0947.

Example 95: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1,3-thiazole-2-carboxamide 4-methylbenzenesulfonate (from 2-bromo-1-(4-fluorophenyl)-ethanone). Yield 16%. HRMS (FAB) calculated for  $C_{17}H_{18}FN_3OS+H_1$  332.1233, found 332.1233.

Example 96: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1,3-thiazole-2-carboxamide 4-methylbenzenesulfonate (from 2-bromo-1-(2-methoxyphenyl)ethanone). Yield 14%. HRMS (FAB) calculated for  $C_{18}H_{21}N_3O_2S+H_1$  344.1432, found 344.1436.

Example 97: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-1,3-thiazole-2-carboxamide 4-methylbenzenesulfonate (from 2-bromo-1-(4-chlorophenyl)-ethanone). Yield 12%. HRMS (FAB) calculated for  $C_{17}H_{18}ClN_3OS+H_1$  348.0937, found 348.0934.

Example 98: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-methyl-1,3-thiazole-2-carboxamide 4-methylbenzenesulfonate (from 1-chloroacetone). Yield 2%. HRMS (FAB) calculated for  $C_{12}H_{17}N_3OS+H_1$  252.1171, found 252.1171.

Example 99: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1,3-thiazole-2-carboxamide 4-methylbenzenesulfonate (from 2-bromo-1-(3-chlorophenyl)-ethanone). Yield 10%. HRMS (FAB) calculated for  $C_{17}H_{18}ClN_3OS+H_1$  348.0937, found 348.0936.

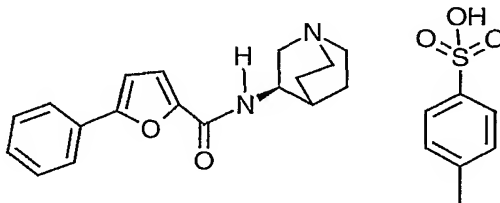
Example 100: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-1,3-thiazole-2-carboxamide 4-methylbenzenesulfonate (from 2-bromo-1-(2-fluorophenyl)-ethanone). Yield 0.4%. MS (ESI) for  $C_{17}H_{18}FN_3OS$  m/z 332.2 (M+H)<sup>+</sup>.

Example 101: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1,3-thiazole-2-carboxamide 4-methylbenzenesulfonate (from 2-bromo-1-(3-fluorophenyl)-ethanone). Yield 6%. MS (ESI) for  $C_{17}H_{18}FN_3OS$  m/z 332.2 (M+H)<sup>+</sup>.

Example 102: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-thien-2-yl-1,3-thiazole-2-carboxamide 4-methylbenzenesulfonate (from 2-bromo-1-thien-2-yl-ethanone). Yield 10%. MS (ESI) for  $C_{15}H_{17}N_3OS_2$  m/z 320.2 (M+H)<sup>+</sup>.

## Example 103

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-furan-2-carboxamide 4-methylbenzenesulfonate:



5 Step 103a: Synthesis of 5-phenyl-furan-2-carbaldehyde.

A solution of 5-bromo-furan-2-carbaldehyde (1.08 g, 6.16 mmol, 1 eq), phenylboronic acid (0.90 g, 7.39 mmol, 1.1 eq), tetrabutylammonium bromide (1.99 g, 6.16 mmol, 1 eq), palladium acetate (30 mg, 0.012 mmol 0.02 eq),  $K_2CO_3$  (2.13 g, 15.4 mmol, 2.5 eq) in water (10 mL) is stirred under nitrogen at room temperature  
 10 overnight. Upon completion, the reaction mixture turns into a black mass floating in the solution. The reaction is diluted with 40 mL water and extracted with EtOAc (3 x 100 mL). The organic layers are combined and stirred with charcoal for 30 min, then dried over  $MgSO_4$  and filtered. The solvent is removed under reduced pressure to give an oil. The product is purified by silica gel chromatography using a Biotage  
 15 Flash 40M column (10% EtOAc/heptane). Yield 70.1%. HRMS (FAB) calculated for  $C_{11}H_8O_2+H$  173.0603, found 173.0607.

Step 103b: Synthesis of 5-phenyl-furan-2-carboxylic acid.

To a solution of the product from Step 103a (0.650 g, 3.78 mmol, 1 eq) in water (5.5 mL), t-BuOH (18.0 mL), and THF (18.0 mL) is added 2-methyl-2-butene  
 20 (3.2 mL, 30.2 mmol, 8 eq), potassium phosphate monobasic (1.54 g, 11.3 mmol, 3 eq), then  $NaClO_2$  (1.03 g, 11.3 mmol, 3 eq) in that order. After four hours, the reaction is complete and diluted with 1 N NaOH (100 mL). The aqueous solution is extracted with ether (2 x 100 mL), and the aqueous layer is acidified with conc. HCl. The resulting solution is extracted with  $CH_2Cl_2$  (3 x 100 mL). The organic layers are  
 25 dried over  $MgSO_4$ , and the solvent removed. The product is purified by silica gel chromatography using a Biotage Flash 40M column (10% EtOAc/1% formic acid/heptane). The solid remaining after removal of the solvent is filtered and

recrystallized from EtOH and water to give the acid as a white crystalline solid (0.499 g, 70.2%). HRMS (FAB) calculated for  $C_{11}H_8O_3+H$  189.0473, found 189.0403.

Step 103c: Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-2-furamide 4-methylbenzenesulfonate.

- 5           The product of Step 103b (0.499 g, 2.65 mmol, 1 eq) and the dihydrochloride salt of R-(+)-aminoquinuclidine (0.528 g, 2.65 mmol, 1 eq) in DIEA (1.38 mL, 7.96 mmol, 3 eq), in THF (17.6 mL) are cooled to 0°C and HATU (1.01 g, 2.65 mmol, 1 eq) is added. The reaction is stirred overnight. The reaction is diluted with  $CH_2Cl_2$  (150 mL). The organic is washed with 1N NaOH, satd  $NaHCO_3$ , and water. The
- 10   organic layer is dried over  $MgSO_4$ , and the solvent is removed under reduced pressure. The product is purified by silica gel chromatography using a Biotage Flash 40M column (8% MeOH/1%  $NH_4OH/CH_2Cl_2$ ). The p-toluene sulfonic acid salt is made, which gave an amorphous solid (1.06 g, 42%). HRMS (FAB) calculated for  $C_{18}H_{20}N_2O_2+H$  297.1603, found 297.1602.

15

#### Examples 104-130

The following compounds are prepared from the requisite boronic acid, furaldehyde, or furan-carboxylic acid according to the procedures for Example 103, making non-critical variations.

- 20   Example 104: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-furan-2-carboxamide toluene sulfonate (from 5-(2-chlorophenyl)-2-furaldehyde). Yield 46%. HRMS (FAB) calculated for  $C_{18}H_{19}ClN_2O_2+H$  331.1213, found 331.1208.
- Example 105: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-furan-2-carboxamide toluene sulfonate (from 5-(3-chlorophenyl)-2-furaldehyde). Yield 72%.
- 25   HRMS (FAB) calculated for  $C_{18}H_{19}ClN_2O_2+H$  331.1213, found 331.1212.
- Example 106: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-furan-2-carboxamide fumarate (from 5-(4-chlorophenyl)-2-furoic acid). Yield 43%. HRMS (FAB) calculated for  $C_{18}H_{19}ClN_2O_2+H$  331.1213, found 331.1216.
- Example 107: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-furan-2-
- 30   carboxamide toluene sulfonate (from 5-(4-bromophenyl)-2-furaldehyde). Yield 38%. HRMS (FAB) calculated for  $C_{18}H_{19}BrN_2O_2+H$  375.0708, found 375.0713.

Example 108: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(trifluoromethyl)-phenyl]-furan-2-carboxamide fumarate (from 5-(2-(trifluoromethyl)-phenyl)-furan-2-carboxylic acid). Yield 78%. HRMS (FAB) calculated for  $C_{19}H_{19}F_3N_2O_2+H_1$  365.1477, found 365.1468.

- 5 Example 109: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(trifluoromethyl)-phenyl]-furan-2-carboxamide fumarate (from 5-(3-(trifluoromethyl)-phenyl)-furan-2-carboxaldehyde). Yield 31%. MS (ESI) for  $C_{19}H_{19}F_3N_2O_2$   $m/z$  365.2 (M+H)<sup>+</sup>.

Example 110: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-furan-2-carboxamide 4-methylbenzenesulfonate (from 5-(2-nitro-phenyl)-furan-2-carboxylic  
10 acid). Yield 82%. HRMS (FAB) calculated for  $C_{18}H_{19}N_3O_4+H_1$  342.1454, found 342.1469.

Example 111: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-furan-2-carboxamide 4-methylbenzenesulfonate (from 5-(3-nitro-phenyl)-furan-2-carboxylic  
15 acid). Yield 66%. HRMS (FAB) calculated for  $C_{18}H_{19}N_3O_4+H_1$  342.1454, found 342.1463.

Example 112: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-furan-2-carboxamide 4-methylbenzenesulfonate (from 5-(4-nitro-phenyl)-furan-2-carboxylic  
acid). Yield 52%. HRMS (FAB) calculated for  $C_{18}H_{19}N_3O_4+H_1$  342.1454, found 342.1465.

- 20 Example 113: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-furan-2-carboxamide 4-methylbenzenesulfonate (from 2-fluorophenylboronic acid). Yield 12%. HRMS (FAB) calculated for  $C_{18}H_{19}FN_2O_2+H_1$  315.1508, found 315.1519.

Example 114: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-furan-2-carboxamide 4-methylbenzenesulfonate (from 3-fluorophenylboronic acid). Yield  
25 29%. HRMS (FAB) calculated for  $C_{18}H_{19}FN_2O_2+H_1$  315.1508, found 315.1519.

Example 115: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-furan-2-carboxamide 4-methylbenzenesulfonate (from 4-fluorophenylboronic acid). Yield 16%. HRMS (FAB) calculated for  $C_{18}H_{19}FN_2O_2+H_1$  315.1508, found 315.1500.

Example 116: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-furan-2-carboxamide hydrochloride (from 2,4-difluorophenylboronic acid). Yield 46%. MS  
30 (ESI) for  $C_{18}H_{18}F_2N_2O_2$   $m/z$  333.2 (M+H)<sup>+</sup>.

Example 117: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,5-difluorophenyl)-furan-2-carboxamide hydrochloride (from 2,5-difluorophenylboronic acid). Yield 33%. MS (ESI) for  $C_{18}H_{18}F_2N_2O_2$   $m/z$  333.2 (M+H)<sup>+</sup>.

Example 118: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-furan-2-carboxamide 4-methylbenzenesulfonate (from 2-methoxyphenylboronic acid). Yield 63%. HRMS (FAB) calculated for  $C_{19}H_{22}N_2O_3+H_1$  327.1708, found 327.1717.

Example 119: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-furan-2-carboxamide 4-methylbenzenesulfonate (from 3-methoxyphenylboronic acid). Yield 83%. HRMS (FAB) calculated for  $C_{19}H_{22}N_2O_3+H_1$  327.1708, found 327.1714.

Example 120: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(trifluoromethoxy)phenyl]-furan-2-carboxamide 4-methylbenzenesulfonate (from 3-trifluoromethoxyphenylboronic acid). Yield 58%. HRMS (FAB) calculated for  $C_{19}H_{19}F_3N_2O_3+H_1$  381.1426, found 381.1440.

Example 121: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-chloro-5-(trifluoromethyl)phenyl]-furan-2-carboxamide fumaric acid (from 2-chloro-5-trifluoromethylphenylboronic acid). Yield 55%. HRMS (FAB) calculated for  $C_{19}H_{18}ClF_3N_2O_2+H_1$  399.1087, found 399.1097.

Example 122: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoro-3-methylphenyl)-furan-2-carboxamide 4-methylbenzenesulfonate (from 4-fluoro-3-methylphenylboronic acid). Yield 77%. HRMS (FAB) calculated for  $C_{19}H_{21}FN_2O_2+H_1$  329.1665, found 329.1661.

Example 123: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-furan-2-carboxamide hydrochloride (from 4-cyanophenylboronic acid). Yield 31%. MS (ESI) for  $C_{18}H_{19}N_3O_2$   $m/z$  322.2 (M+H)<sup>+</sup>.

Example 124: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-thien-2-yl-furan-2-carboxamide 4-methylbenzenesulfonate (from 2-thiopheneboronic acid). MS (ESI) for  $C_{16}H_{18}N_2O_2S$   $m/z$  303.2 (M+H)<sup>+</sup>.

Example 125: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-thien-3-yl-furan-2-carboxamide 4-methylbenzenesulfonate (from 3-thiopheneboronic acid). MS (ESI) for  $C_{16}H_{18}N_2O_2S$   $m/z$  303.2 (M+H)<sup>+</sup>.

Example 126: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-furan-2-carboxamide 4-methylbenzenesulfonate (from 5-bromo-furan-2-carboxylic acid). Yield 75%. MS (ESI) for  $C_{12}H_{15}N_2O_2Br$   $m/z$  299.0 (M+H)<sup>+</sup>.

Example 127: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-nitro-furan-2-carboxamide hydrochloride (from 5-nitro-furan-2-carboxylic acid). Yield 63%. MS (ESI) for  $C_{12}H_{15}N_3O_4$   $m/z$  265.1 (M+H)<sup>+</sup>.

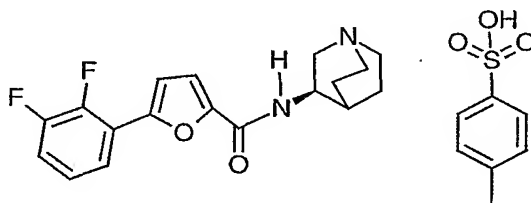
Example 128: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4,5-dimethyl-furan-2-carboxamide hydrochloride (from 4,5-dimethyl-furan-2-carboxylic acid). Yield 75%. HRMS (FAB) calculated for  $C_{14}H_{20}N_2O_2+H$  249.1603, found 249.1593.

Example 129: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloro-2-nitrophenyl)-furan-2-carboxamide hydrochloride (from 4-chloro-2-nitrophenyl)-furan-2-carboxylic acid). Yield 7%. HRMS (FAB) calculated for  $C_{18}H_{18}ClN_3O_4+H$  376.1064, found 376.1067.

Example 130: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methyl-2-nitrophenyl)-furan-2-carboxamide hydrochloride (from 4-methyl-2-nitrophenyl)-furan-2-carboxylic acid). Yield 56%. MS (ESI) for  $C_{19}H_{21}N_3O_4$   $m/z$  356.2 (M+H)<sup>+</sup>.

### Example 131

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3-difluorophenyl)-furan-2-carboxamide 4-methylbenzenesulfonate:



20

Step 131a: A solution of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-furan-2-carboxamide (0.258 g, 0.85 mmol, 1 eq), 2,3-difluorophenylboronic acid (0.147 g, 0.93 mmol, 1.1 eq), tetrabutylammonium bromide (0.244 g, 0.85 mmol, 1 eq), palladium acetate (3.8 mg, 0.017 mmol, 0.02 eq),  $K_2CO_3$  (0.41 g, 2.97 mmol, 3.5 mmol) and water 1.4 mL is stirred under argon overnight. The reaction forms a brownish insoluble lump, but is complete by HPLC. The reaction is purified by silica gel chromatography using a Biotage Flash 40 M column (10%MeOH/1%TEA/ $CH_2Cl_2$ ). The p-toluenesulfonic acid salt of 5-(2,3-difluoro-phenyl)-furan-2-carboxylic acid (1-

aza-bicyclo[2.2.2]oct-3-yl)-amide is synthesized and recrystallized from IPA/MeOH/ether, and the product is recovered as a crystalline solid (0.135 g, 31%). HRMS (FAB) calculated for  $C_{18}H_{18}F_2N_2O_2+H_1$  333.1414, found 333.1418.

#### Examples 132-146

5           The following compounds are prepared from the requisite boronic acid and N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-furan-2-carboxamide as stated for Example 131, making non-critical variations.

Example 132: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenyl)-furan-2-carboxamide hydrochloride (from 3,4-difluorophenyl-boronic acid). Yield 24%.

10   HRMS (FAB) calculated for  $C_{18}H_{18}F_2N_2O_2+H_1$  333.1414, found 333.1418.

Example 133: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-furan-2-carboxamide hydrochloride (from 3,5-difluorophenyl-boronic acid). Yield 60%.

HRMS (FAB) calculated for  $C_{18}H_{18}F_2N_2O_2+H_1$  333.1414, found 333.1424.

Example 134: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxy-phenyl)-furan-2-  
15   carboxamide hydrochloride (from 4-methoxy-phenyl-boronic acid). Yield 17%.

HRMS (FAB) calculated for  $C_{19}H_{22}N_2O_3+H_1$  327.1708, found 327.1707.

Example 135: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-furan-2-carboxamide 4-methylbenzenesulfonate (from *o*-tolylboronic acid). Yield 26%.

HRMS (FAB) calculated for  $C_{19}H_{22}N_2O_2+H_1$  311.1759, found 311.1763.

20   Example 136: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-furan-2-carboxamide 4-methylbenzenesulfonate (from *m*-tolylboronic acid). HRMS (FAB) calculated for  $C_{19}H_{22}N_2O_2+H_1$  311.1759, found 311.1752.

Example 137: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-furan-2-carboxamide 4-methylbenzenesulfonate (from *p*-tolylboronic acid). Yield 10 %.

25   HRMS (FAB) calculated for  $C_{19}H_{22}N_2O_2+H_1$  311.1759, found 311.1752.

Example 138: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(trifluoromethoxy)phenyl]-furan-2-carboxamide 4-methylbenzenesulfonate (from 2-(trifluoromethoxy)phenyl-boronic acid). Yield 19%. MS (ESI) for  $C_{19}H_{19}F_3N_2O_3$   $m/z$  381.3 ( $M+H$ )<sup>+</sup>.

Example 139: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(trifluoromethoxy)phenyl]-  
30   furan-2-carboxamide 4-methylbenzenesulfonate (from 4-trifluoromethoxyphenyl-

boronic acid). Yield 61%. HRMS (FAB) calculated for  $C_{19}H_{19}F_3N_2O_3+H_1$  381.1426, found 381.1434.

Example 140: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-tert-butylphenyl)-furan-2-carboxamide 4-methylbenzenesulfonate (from 4-*tert*-butylphenyl-boronic acid). Yield 5 73%. HRMS (FAB) calculated for  $C_{22}H_{28}N_2O_2+H_1$  353.2229, found 353.2218.

Example 141: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(1-benzothien-2-yl)-furan-2-carboxamide 4-methylbenzenesulfonate (from benzothiophene-2-boronic acid). Yield 17%. HRMS (FAB) calculated for  $C_{20}H_{20}N_2O_2S+H_1$  353.1324, found 353.1326.

Example 142: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-quinolin-3-yl-furan-2-10 carboxamide 4-methylbenzenesulfonate (from 3-quinoline-boronic acid). Yield 9%. MS (ESI) for  $C_{21}H_{21}N_3O_2$   $m/z$  348.3 (M+H)<sup>+</sup>.

Example 143: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-ethylphenyl)-furan-2-carboxamide hydrochloride (from 4-ethylphenyl-boronic acid). MS (ESI) for  $C_{20}H_{24}N_2O_2$   $m/z$  325.3 (M+H)<sup>+</sup>.

15 Example 144: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-isopropylphenyl)-furan-2-carboxamide hydrochloride (from 4-isopropylphenyl-boronic acid). MS (ESI) for  $C_{21}H_{26}N_2O_2$   $m/z$  339.3 (M+H)<sup>+</sup>.

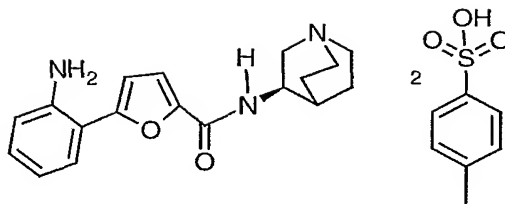
Example 145: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoro-4-methoxyphenyl)-furan-2-carboxamide hydrochloride (from 3-fluoro-4-methoxy-boronic acid). MS 20 (ESI) for  $C_{19}H_{21}FN_2O_2$   $m/z$  345.2 (M+H)<sup>+</sup>.

Example 146: N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(1-benzofuran-2-yl)-furan-2-carboxamide 4-methylbenzenesulfonate (from benzofuran-2-boronic acid). MS (ESI) for  $C_{20}H_{20}N_2O_3$   $m/z$  337.2 (M+H)<sup>+</sup>.

25

## Example 147

5-(2-Aminophenyl)-N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-furan-2-carboxamide bis(4-methylbenzenesulfonate):



Step 147a: To a solution of N-[(3R)-1-aza-bicyclo[2.2.2]oct-3-yl]-5-(2-nitro-phenyl)-furan-2-carboxamide (1.65 g, 3.22 mmol, 1 eq) in 100 mL EtOH was added Pd/C (50 mg). This mixture was placed on a Parr shaker under 40 psi hydrogen  
 5 overnight. The palladium is removed by filtration over a pad of celite, and the solvent is removed. The p-toluenesulfonic acid salt of 5-(2-amino-phenyl)-furan-2-carboxylic acid (1-aza-bicyclo[2.2.2]oct-3-yl)-amide is purified by recrystallization from IPA/MeOH/ether to give a crystalline solid (0.783 g, 50%). HRMS (FAB) calculated for C<sub>18</sub>H<sub>21</sub>N<sub>3</sub>O<sub>2</sub>+H<sub>1</sub> 312.1712, found 312.1717.

10

#### Examples 148-149

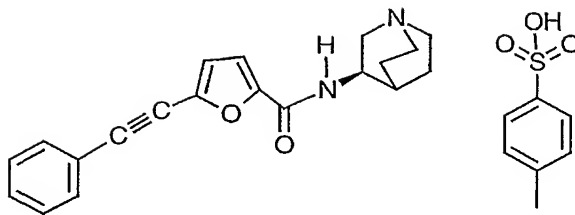
The following compounds are prepared from the requisite nitro amide according to the procedures for Example 147, making non-critical variations..

Example 148: 5-(4-aminophenyl)-N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-furan-2-  
 15 carboxamide tris(4-methylbenzenesulfonate) (from N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-2-furamide). Yield 84%. HRMS (FAB) calculated for C<sub>18</sub>H<sub>21</sub>N<sub>3</sub>O<sub>2</sub>+H<sub>1</sub> 312.1712, found 312.1727.

Example 149: 5-(2-amino-4-methylphenyl)-N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-  
 20 furan-2-carboxamide dihydrochloride (from 5-(4-methyl-2-nitrophenyl)-N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-furamide dihydrochloride). Yield 55%. HRMS (FAB) calculated for C<sub>19</sub>H<sub>23</sub>N<sub>3</sub>O<sub>2</sub>+H 326.1868, found 326.1871.

#### Example 150

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylethynyl)-furan-2-carboxamide  
 25 4-methylbenzenesulfonate:



Step 150a: Preparation of 5-phenylethynyl-furan-2-aldehyde.

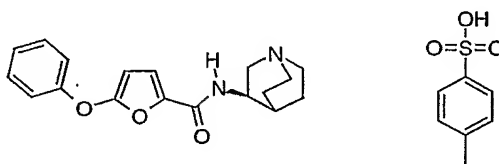
To a solution of 5-bromofuraldehyde (1.0 g, 5.71 mmol, 1eq), copper (I) iodide (0.163 g, 0.857 mmol, 0.15 eq), trans-dichlorobis(triphenylphosphine) palladium(II) (0.20 g, 0.287 mmol, 0.05 eq), and TEA (3.98 mL, 28.6 mmol, 5 eq) in THF (45 mL) is added dropwise phenyl acetylene (1.25 mL, 11.4 mmol, 2 eq). After 48 hours, the reaction appears complete. The reaction is filtered over a pad of celite, and the solvent is removed under reduced pressure. The reaction is purified by silica gel chromatography using a Biotage Flash 40M column (10% EtOAc/heptane) to give a yellow orange crystalline solid (0.765 g, 68.3%). MS (ESI) for C<sub>13</sub>H<sub>8</sub>O<sub>2</sub> *m/z* 197.1 (M+H)<sup>+</sup>.

Step 150b: Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylethynyl)-furan-2-carboxamide 4-methylbenzenesulfonate.

Following the general procedure of Example 103, making non-critical variations but starting with 5-phenylethynyl-furan-2-aldehyde, N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylethynyl)-furan-2-carboxamide 4-methylbenzenesulfonate is synthesized and recovered as a crystalline solid (0.692 g, 74%). HRMS (FAB) calculated for C<sub>20</sub>H<sub>20</sub>N<sub>2</sub>O<sub>2</sub>+H<sub>1</sub> 321.1603, found 321.1595.

## Example 151

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-furan-2-carboxamide 4-methylbenzenesulfonate:



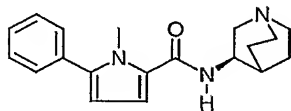
Step 151a: Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-furan-2-carboxamide 4-methylbenzenesulfonate.

A solution of N-[(3R)-1-aza-bicyclo[2.2.2]oct-3-yl]-5-bromo-furan-2-carboxamide (0.200 g, 0.42 mmol, 1 eq), sodium phenoxide (0.500 g, 4.3 mmol, 10.1 eq), in DMSO (5 mL) is stirred under nitrogen at room temperature overnight. The reaction is diluted with 25 mL water and extracted with CH<sub>2</sub>Cl<sub>2</sub> (50 mL). The organic layer is washed with water (3x25 mL), satd NaHCO<sub>3</sub>, brine, and dried over MgSO<sub>4</sub>. The p-toluenesulfonic acid salt is prepared and recrystallized from IPA/MeOH/ether, and the product is recovered as a crystalline solid (70%). HRMS (FAB) calculated for C<sub>18</sub>H<sub>20</sub>N<sub>2</sub>O<sub>3</sub>+H 313.1552, found 313.1558.

10

## Example 152

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-1-methyl-5-phenyl-1H-pyrrole-2-carboxamide:



Step 152a: Preparation of methyl 5-bromo-1-methyl-1H-pyrrole-2-carboxylate.

15

To a dry flask is added methyl 1-methyl-1H-pyrrole-2-carboxylate (12.0 g, 86.4mmol) and 150 mL of dry CH<sub>2</sub>Cl<sub>2</sub>, and the flask is wrapped in foil and purged with nitrogen. N-bromosuccinimide (16.2 g, 90.7 mmol) is added in one portion and the mixture is stirred at room temperature for 0.5 h. The reaction mixture is washed with water (50 mL) and brine (50 mL), dried over MgSO<sub>4</sub>, filtered, and concentrated under reduced pressure. Fractional distillation gives 12.0 g of methyl 5-bromo-1-methyl-1H-pyrrole-2-carboxylate as a yellow oil (64% yield). MS for C<sub>7</sub>H<sub>8</sub>NO<sub>2</sub>Br (ESI) (M)<sup>+</sup> m/z 217.1.

20

Step 152b: Preparation of methyl 1-methyl-5-phenyl-1H-pyrrole-2-carboxylate.

25

The product from step 152a is added to a solution of tetrakis(triphenylphosphine)palladium(0) (0.530 g, 0.459mmol) in 90 mL of ethylene glycol dimethyl ether. The resulting solution is stirred under nitrogen for 5 min and then phenylboronic acid (1.34 g, 11.0 mmol) is added followed by a solution of Na<sub>2</sub>CO<sub>3</sub> (19.5 g, 183 mmol) in 90 mL of H<sub>2</sub>O. The mixture is heated at reflux for 24 hours. The reaction mixture is allowed to cool to rt, 100 mL of CH<sub>2</sub>Cl<sub>2</sub> is added, and the layers are separated. The aqueous layer is extracted with CH<sub>2</sub>Cl<sub>2</sub> (3 x 50 mL) and

30

combined organic layers are dried over  $\text{MgSO}_4$ , filtered, and concentrated *in vacuo*. The crude product is purified by flash column chromatography (5 % EtOAc in hexane) to give 1.89 g of methyl 1-methyl-5-phenyl-1H-pyrrole-2-carboxylate as a yellow oil (96% yield). MS for  $\text{C}_{13}\text{H}_{13}\text{NO}_2$  (ESI)  $(\text{M}+\text{H})^+$   $m/z$  216.1.

5 Step 152c: Preparation of 1-methyl-5-phenyl-1H-pyrrole-2-carboxylic acid.

Lithium hydroxide (1.39 g, 33.2 mmol) is added to a solution of the product from Step 152b (1.43 g, 6.64 mmol) in 96 mL of a 1.25:1:1  $\text{H}_2\text{O}:\text{MeOH}:\text{THF}$  solvent mixture. The reaction is stirred at  $50^\circ\text{C}$  for 2 h. Aqueous HCl (1N, 50 mL) is added and the resultant precipitate is collected by filtration, washed with water, and dried to  
10 give 0.851 g of 1-methyl-5-phenyl-1H-pyrrole-2-carboxylic acid as a tan solid (64% yield). MS for  $\text{C}_{12}\text{H}_{11}\text{NO}_2$  (ESI)  $(\text{M}-\text{H})^+$   $m/z$  200.1.

Step 152d: Preparation of *N*-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-1-methyl-5-phenyl-1H-pyrrole-2-carboxamide.

To a solution of (*R*)-(+)-3-aminoquinuclidine dihydrochloride (0.831 g, 4.17  
15 mmol) in 55 mL of dry THF is added DIEA (2.08 mL, 11.9 mmol). After allowing the mixture to stir under nitrogen for 15 min, the product from Step 152c (0.800 g, 3.98 mmol) is added. The mixture is allowed to stir for another 15 min, cooled in an ice bath, and HATU (1.59 g, 4.17 mmol) is added. The reaction mixture is stirred on an ice bath for 0.5 h, then at rt for an additional 2 h. The mixture is diluted with 60  
20 mL  $\text{CH}_2\text{Cl}_2$ , washed with 50 mL of 1N NaOH and 50 mL of satd  $\text{NaHCO}_3$  solution, dried over  $\text{MgSO}_4$ , filtered, and concentrated *in vacuo*. The crude product is purified by flash column chromatography (10 % EtOAc, 1 %  $\text{NH}_4\text{OH}$  in  $\text{CH}_2\text{Cl}_2$ ) to give 0.383 grams of *N*-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-1-methyl-5-phenyl-1H-pyrrole-2-carboxamide (31%) as a white solid. MS for  $\text{C}_{19}\text{H}_{23}\text{N}_3\text{O}$  (ESI)  $(\text{M}+\text{H})^+$   $m/z$  310.3.

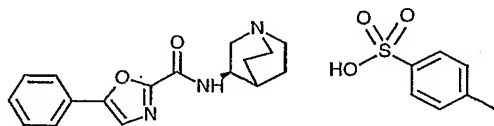
25

### Examples 153-157

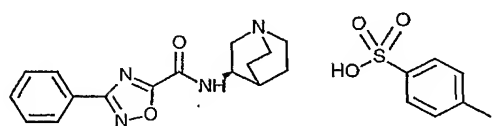
These examples are prepared using the coupling procedure for Example 103c, making non-critical variations and using the appropriate carboxylic acids.

Example 153: *N*-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1,3-oxazole-2-carboxamide 4-methylbenzenesulfonate (from 5-phenyl-1,3-oxazole-2-carboxylic  
30

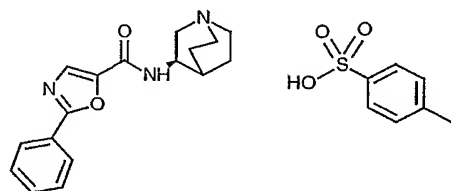
acid, see Saito, S.; Tanaka, C. *J. Pharm. Sci. Japan* 76, 1956, 305-7). Yield 77%.  
 HRMS (FAB) calculated for  $C_{17}H_{19}N_3O_2+H_1$  298.1555, found 298.1558.



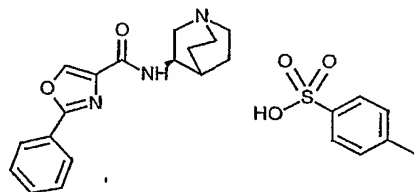
Example 154: *N*-[(3*R*)-1-azabicyclo[2.2.2]oct-3-yl]-3-phenyl-1,2,4-oxadiazole-5-  
 5 carboxamide 4-methylbenzenesulfonate (from 3-phenyl-1,2,4-oxadiazole-5-carboxylic  
 acid, see Wurm. *Chem. Ber.*; 22; 1889; 3133). Yield 54%. HRMS (FAB) calculated  
 for  $C_{16}H_{18}N_4O_2+H_1$  299.1508, found 299.1512.



Example 155: *N*-[(3*R*)-1-azabicyclo[2.2.2]oct-3-yl]-2-phenyl-1,3-oxazole-5-  
 10 carboxamide 4-methylbenzenesulfonate (from 2-phenyl-1,3-oxazole-5-carboxylic  
 acid, see Belen'kii, L. I.; Cheskis, M. A.; Zvolinskii, V. P.; Obukhov, A. E. *Chem.*  
*Heterocycl. Compd. (Engl. Transl.)*; 22; 1986; 654-663). Yield 16%. HRMS (FAB)  
 calculated for  $C_{17}H_{19}N_3O_2+H_1$  298.1555, found 298.1555.

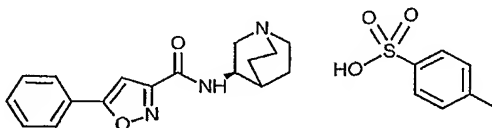


Example 156: *N*-[(3*R*)-1-azabicyclo[2.2.2]oct-3-yl]-2-phenyl-1,3-oxazole-4-  
 15 carboxamide 4-methylbenzenesulfonate (from 2-phenyl-1,3-oxazole-4-carboxylic  
 acid, see Korte, F.; Stoeriko, K. *Chem. Ber.*; 93; 1960; 1033-1042). Yield 22%.  
 HRMS (FAB) calculated for  $C_{17}H_{19}N_3O_2+H_1$  298.1555, found 298.1559.



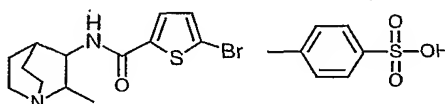
Example 157: *N*-[(3*R*)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenylisoxazole-3-  
 20 carboxamide 4-methylbenzenesulfonate (from 5-phenylisoxazole-3-carboxylic acid,

see Vaughan, W.R.; Spencer, J.L. *J.Org.Chem.*; 25; 1960; 1160-1164). Yield 76%.  
 HRMS (FAB) calculated for  $C_{17}H_{19}N_3O_2 + H_1$  298.1555, found 298.1556.



Example 158

- 5            5-Bromo-N-(2-methyl-1-azabicyclo[2.2.2]oct-3-yl)thiophene-2-carboxamide  
 4-methylbenzenesulfonate:



Step 158a: Preparation of 2-methylenequinuclidin-3-one.

A mixture of 2-methylene-3-quinuclidinone dihydrate hydrochloride (25.7g,  
 10 0.1225mol, 1eq) and  $K_2CO_3$  (67.0g, 0.4848mol, 4eq) is dissolved in 125mL water and  
 200mL  $CH_2Cl_2$  and is stirred vigorously. After 16h, the layers are separated and the  
 aqueous layer is extracted with  $CH_2Cl_2$ . The combined organic layers are dried over  
 $MgSO_4$ , filtered and concentrated to give 14.75g (88%) of 2-methylenequinuclidin-3-  
 one as a yellow oil. MS (ESI) for  $C_8H_{11}NO$   $m/z$  138.1 ( $M+H$ )<sup>+</sup>.

- 15 Step 158b: Preparation of 2-methylquinuclidin-3-one hydrochloride.

The product from Step 158a (14.75g, 0.1075mol, 1eq), formic acid (10.4g,  
 0.2150mol, 2eq) and  $(Ph_3P)_3RuCl_2$  (0.21g, 0.21mmol) are dissolved in 100mL THF.  
 The mixture is heated under reflux. Fresh portions of catalyst (0.58g, 0.59mmol,  
 (total)) and formic acid (1.2g, 0.026mol) are added periodically over the course of the  
 20 reaction. After 72h, the mixture is concentrated *in vacuo*. The residue is taken up in  
 ether and excess HCl in dioxane (27mL, 4.0M) is added. The solids are washed with  
 ether and recrystallized from EtOH to afford 14.4g (76%) of 2-methylquinuclidin-3-  
 one hydrochloride as a white solid. MS (ESI) for  $C_8H_{13}NO$   $m/z$  140.2 ( $M+H$ )<sup>+</sup>.

Step 158c: Preparation of (3E/Z)-2-methyl-1-azabicyclo[2.2.2]octan-3-one oxime.

- 25            The product from Step 158b (3.2g, 23.0mmol, 1eq) and hydroxylamine  
 hydrochloride (1.6g, 23.0mmol, 1eq) are dissolved in 20mL EtOH/pyridine (4:1) and  
 stirred at room temperature. After 5 days, water and solid NaOH are added to adjust

pH to pH 11. The mixture is extracted with several portions of  $\text{CHCl}_3$ . The combined organic layers are dried over  $\text{MgSO}_4$ , filtered and concentrated to give 3.42g (96%) of 2-methyl-1-azabicyclo[2.2.2]octan-3-one oxime as a 1:2.6 mixture of oxime isomers. Partial  $^1\text{H}$  NMR ( $\text{CDCl}_3$ , 300 MHz)  $\delta$  1.53ppm (d, 2- $\text{CH}_3$ , 0.8H), 1.38 (d, 2- $\text{CH}_3$ , 2.2H). MS (ESI) for  $\text{C}_8\text{H}_{14}\text{N}_2\text{O}$   $m/z$  154.8 ( $\text{M}+\text{H}$ ) $^+$ .

Step 158d: Preparation of 2-methylquinuclidin-3-amine dihydrochloride

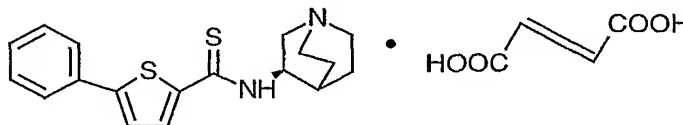
Sodium (7.0g, 0.303mol, 10eq) is added in portions to a solution of 2-methyl-1-azabicyclo[2.2.2]octan-3-one oxime from Step 158c (4.65g, 30.2mmol, 1eq) in 100 mL *n*-propanol. The mixture is heated under reflux. After about 12h, the mixture is cooled and 80mL of water is added. The layers are separated, and the aqueous layer is extracted with  $\text{CHCl}_3$ . The combined organic layers are dried over  $\text{MgSO}_4$  and filtered. An excess of HCl in dioxane (15mL, 4.0M) is added to the solution and the solvent is removed to give 6.0g (93%) of 2-methylquinuclidin-3-amine dihydrochloride as an oil. A hygroscopic solid was obtained by trituration of the oil in hot IPA. MS (ESI) for  $\text{C}_8\text{H}_{16}\text{N}_2$   $m/z$  141.3 ( $\text{M}+\text{H}$ ) $^+$ .

Step 158e: Preparation of 5-bromo-N-(2-methyl-1-azabicyclo[2.2.2]oct-3-yl)thiophene-2-carboxamide 4-methylbenzenesulfonate.

5-Bromothiophene-2-carboxylic acid (1.63g, 7.88mmol, 1eq), the product from Step 158d (1.68g, 7.88mmol, 1eq) and HATU (2.97g, 7.81mmol, 1eq) are suspended in 60mL  $\text{CH}_3\text{CN}$ . The mixture is cooled in an ice bath and DIEA (8.2mL, 47.28mmol, 6eq) is added dropwise. The mixture is allowed to warm to room temperature and stirred overnight. EtOAc and satd  $\text{NaHCO}_3$  are added. The aqueous layer is extracted with EtOAc. The combined organic layers are washed with 1N NaOH, dried over  $\text{MgSO}_4$ , filtered and concentrated. The residue is purified by silica gel chromatography using a Biotage Flash 40S column (90:9:1  $\text{CHCl}_3/\text{MeOH}/\text{NH}_4\text{OH}$ ) to provide 1.26g (49%) of 5-bromo-N-(2-methyl-1-azabicyclo[2.2.2]oct-3-yl)thiophene-2-carboxamide as a trans/cis mixture of isomers. The *p*-toluenesulfonic acid salt of the product is made and recrystallized from IPA/ether to give the product as light yellow solid. MS (ESI) for  $\text{C}_{13}\text{H}_{17}\text{BrN}_2\text{OS}$   $m/z$  328.9/330.9 ( $\text{M}+\text{H}$ ) $^+$ . Reverse phase HPLC (ZORBAX Eclipse XDB-C8, 4.6mm x 15cm, 80:12:8  $\text{H}_2\text{O}/\text{CH}_3\text{CN}/\text{IPA}$ ) revealed a 95:5 trans/cis mixture of isomers.

## Example 159

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-thiophene-2-carbothioamide fumarate:



5 Step 159a: Preparation of methyl 5-phenyl-thiophene-2-carbodithioate.

To a cooled (-10 to 0°C) solution of n-BuLi (22.7 mL, 33.4 mmol) in THF (10 mL) is added dropwise a solution of 2-phenyl-thiophene (5.46 g, 34.0 mmol) in THF (15 mL). The resulting green solution is stirred at 0°C. After 30 minutes, a solution of copper (I) bromide (0.87 g, 6.1 mmol) and lithium bromide (1.29 g, 14.9 mmol) in THF (20 mL) is added to the cooled reaction solution over several minutes. The resulting dark green solution is stirred at 0°C for 15 minutes, at which time, carbon disulfide (2.0 mL, 34.0 mmol) is added dropwise over 15 minutes. The resulting dark brown solution is stirred for 30 minutes, then iodomethane (2.9 mL, 46.4 mmol) is added dropwise to the reaction solution over 5 minutes. The resulting dark brown solution is allowed to warm to room temperature and stirred for 1 hour, then is quenched with a solution of potassium cyanide in water (100 mL). The biphasic mixture was diluted with EtOAc and washed with brine, dried over MgSO<sub>4</sub>, filtered, and concentrated to give a dark orange solid (8.5 grams) which is purified with flash chromatography on silica gel (eluent: gradient of heptane to 2% THF/heptane) to give an orange solid. Yield 34%. HRMS (FAB) calculated for C<sub>12</sub>H<sub>10</sub>S<sub>3</sub>+H 251.0023, found 251.0023.

Step 159b: Preparation of N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-thiophene-2-carbothioamide fumarate.

25 A solution of the product of Step 159a (0.49 g, 2.0 mmol) and (R)-3-aminoquinulidine (0.55 g, 4.4 mmol) in THF is stirred at 50°C for 19 hours. The crude reaction mixture is absorbed on to silica gel and purified by flash chromatography (gradient of 7% [9:1 MeOH/NH<sub>4</sub>OH]/CH<sub>2</sub>Cl<sub>2</sub> to 9% [9:1 MeOH/NH<sub>4</sub>OH]/CH<sub>2</sub>Cl<sub>2</sub>). The fumarate salt is prepared and crystallized. Yield 75%. HRMS (FAB) calculated for C<sub>18</sub>H<sub>20</sub>N<sub>2</sub>S<sub>2</sub>+H 329.1146, found 329.1151.

The present invention also includes, by representation but not limitation, any one of the following or combination of the following compounds and pharmaceutically acceptable salts thereof, both of which can be made by one of ordinary skill in the art using the procedures provided making non-critical changes:

5 N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3'-bithiophene-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-benzyloxyphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-benzyloxyphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoro-4-benzyloxyphenyl)-thiophene-2-carboxamide; 5-(2-aminophenyl)-N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-pyridin-3-yl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5'-methyl-2,2'-bithiophene-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5'-chloro-2,2'-bithiophene-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-nitro-thiophene-2-carboxamide; 5-(aminomethyl)-N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-cyano-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-methoxy-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-[2,2']bithiophenyl-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(methylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-acetyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-methyl-thiophene-2-

carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(acetylamino)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-trifluoromethyl-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methyl-5-trifluoromethyl-2H-pyrazole-3-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylthiazol-4-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(3-chlorophenyl)-vinyl]-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl-sulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl-sulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chloro-4-fluoro-phenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3-dichlorophenyl-sulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4,5-trichlorophenyl-sulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl-sulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoro-4-hydroxyphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(phenylsulfanyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[(4-chlorophenyl)-sulfanyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-phenoxy-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[(4-fluorophenyl)-sulfanyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(methylsulfanyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-chlorophenoxy)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-phenyl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2,4-

dimethyl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-fluorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-fluorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-fluorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-hydroxyphenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-methylphenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[4-(benzyloxy)phenyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-pyridin-3-yl-1,3-thiazole-4-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-2-phenyl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-chlorophenyl)-4-methyl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-2-pyridin-2-yl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-2-pyridin-4-yl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(methylamino)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylsulfanyl)-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-chlorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenoxy)-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-fluorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-chlorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-fluorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-chlorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1H-1,2,4-triazole-3-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1,3,4-oxadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(methylthio)-1,3,4-oxadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-

methoxyphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-5-phenyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-methyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-thien-2-yl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-

bromophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(trifluoromethyl)-phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(trifluoromethyl)-phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,5-difluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(trifluoromethoxy)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-chloro-5-(trifluoromethyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoro-3-methylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-thien-2-yl-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-thien-3-yl-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-nitro-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4,5-dimethyl-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloro-2-nitrophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methyl-2-nitrophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3-difluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-

(2-methylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(trifluoromethoxy)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(trifluoromethoxy)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-tert-butylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(1-benzothien-2-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-quinolin-3-yl-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-ethylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-isopropylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoro-4-methoxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(1-benzofuran-2-yl)-furan-2-carboxamide; 5-(2-aminophenyl)-N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-furan-2-carboxamide; 5-(4-aminophenyl)-N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-furan-2-carboxamide; 5-(2-amino-4-methylphenyl)-N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylethynyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-1-methyl-5-phenyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-3-phenyl-1,2,4-oxadiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-phenyl-1,3-oxazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-phenyl-1,3-oxazole-4-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenylisoxazole-3-carboxamide; 5-bromo-N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-thiophene-2-carbothioamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetamidophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetamidophenyl)-

thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetamidophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-difluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-difluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-difluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-carbamoylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-carbamoylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-carbamoylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-sulfamoylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-sulfamoylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-sulfamoylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-

azabicyclo[2.2.2]oct-3-yl]-5-(2-methanesulfonylaminophenylsulfanyl)-thiophene-2-  
 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
 methanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(4-methanesulfonylaminophenylsulfanyl)-thiophene-2-  
 5 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
 trifluoromethanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethanesulfonylaminophenylsulfanyl)-  
 thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
 trifluoromethanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-  
 10 azabicyclo[2.2.2]oct-3-yl]-5-(2-difluoroacetamidophenylsulfanyl)-thiophene-2-  
 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
 difluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(4-difluoroacetamidophenylsulfanyl)-thiophene-2-  
 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-carbamoylphenylsulfanyl)-  
 15 thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
 carbamoylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-  
 3-yl]-5-(4-carbamoylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(2-sulfamoylphenylsulfanyl)-thiophene-2-carboxamide;  
 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-sulfamoylphenylsulfanyl)-thiophene-2-  
 20 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-sulfamoylphenylsulfanyl)-  
 thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
 hydroxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
 yl]-5-(3-hydroxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenylsulfanyl)-thiophene-2-carboxamide; N-  
 25 [(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenylsulfanyl)-thiophene-2-  
 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenylsulfanyl)-  
 thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
 aminophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
 yl]-5-(2-acetamidophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-  
 30 azabicyclo[2.2.2]oct-3-yl]-5-(3-acetamidophenoxy)-thiophene-2-carboxamide; N-  
 [(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetamidophenoxy)-thiophene-2-  
 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoroacetamidophenoxy)-  
 thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-

trifluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-difluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-difluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-difluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-carbamoylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-carbamoylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-carbamoylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-sulfamoylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-sulfamoylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-sulfamoylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenoxy)-thiophene-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenoxy)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenoxy)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-thiophene-2-carboxamide;  
5 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethylphenyl)-thiophene-2-  
10 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethylphenyl)-  
thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
trifluoromethylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(2-acetylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(3-acetylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
15 yl]-5-(4-acetylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(2-cyanophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(3-cyanophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(4-cyanophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(2-morpholin-4-yl-phenyl)-thiophene-2-carboxamide; N-[(3R)-1-  
20 azabicyclo[2.2.2]oct-3-yl]-5-(3-morpholin-4-yl-phenyl)-thiophene-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-morpholin-4-yl-phenyl)-thiophene-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenoxy)-  
thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
methoxyphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-  
25 (4-methoxyphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(2-methylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-  
3-yl]-5-(3-methylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenoxy)-thiophene-2-carboxamide; N-[(3R)-  
1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethylphenoxy)-thiophene-2-carboxamide;  
30 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethylphenoxy)-thiophene-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethylphenoxy)-  
thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
acetylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-

(3-acetylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-  
5 (4-cyanophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-morpholin-4-yl-phenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-morpholin-4-yl-phenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-morpholin-4-yl-phenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenylsulfanyl)-  
10 thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenylsulfanyl)-thiophene-2-  
15 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
20 trifluoromethylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
25 cyanophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-morpholin-4-yl-phenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-morpholin-4-yl-  
30 phenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-morpholin-4-yl-phenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylpyridin-4-yl)-thiophene-2-carboxamide; N-

[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methylpyridin-3-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylpyridin-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxypyridin-4-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methoxypyridin-3-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxypyridin-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-3-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-3-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylpyridin-4-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methylpyridin-3-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylpyridin-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxypyridin-4-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methoxypyridin-3-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxypyridin-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-3-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-2-yloxy)-

thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylpyridin-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methylpyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylpyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxypyridin-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methoxypyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxypyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiophen-4-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiophen-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiophen-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiophen-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(furan-4-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylfuran-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyfuran-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorofuran-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyloxazol-2-yl)-

thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyoxazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorooxazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-oxazol-5-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methyloxazol-5-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyoxazol-5-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorooxazol-5-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-5-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylthiazol-5-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxythiazol-5-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorothiazol-5-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylthiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxythiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorothiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methyloxazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyoxazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorooxazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]thiadiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]thiadiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]thiadiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]thiadiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]oxadiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]oxadiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]oxadiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]oxadiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-

azabicyclo[2.2.2]oct-3-yl]-5-(thiophen-4-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiophen-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiophen-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiophen-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(furan-4-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylfuran-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyfuran-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorofuran-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyloxazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyoxazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorooxazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-5-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methyloxazol-5-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyoxazol-5-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorooxazol-5-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-5-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylthiazol-5-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxythiazol-5-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorothiazol-5-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylthiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxythiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorothiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methyloxazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyoxazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-

3-yl]-5-(4-chlorooxazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]thiadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]thiadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]thiadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]thiadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]oxadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]oxadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]oxadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]oxadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiophen-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiophen-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiophen-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiophen-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(furan-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylfuran-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyfuran-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorofuran-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyloxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyoxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorooxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methyloxazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyoxazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorooxazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiazol-2-ylsulfanyl)-thiophene-2-

carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylthiazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxythiazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorothiazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylthiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxythiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorothiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methyloxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyoxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorooxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]thiadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]thiadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]thiadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]thiadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]oxadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]oxadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]oxadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]oxadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrrole-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(isothiazol-3-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(isoxazol-3-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3H-imidazol-4-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-(4-hydroxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-(3-acetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-(2-methanesulfonylaminophenyl)-thiophene-

2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-(pyridin-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-(2-methylpyridin-4-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-(4-trifluoromethylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-(2-acetylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-phenyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-5-phenyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-cyano-5-phenyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-methoxy-5-phenyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-phenyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-methyl-4-phenyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-cyano-4-phenyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-methoxy-4-phenyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-bromo-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-methylsulfanyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-methyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-cyano-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-bromo-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-methylsulfanyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-methyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-cyano-thiophene-2-carboxamide;

N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methanesulfonylaminophenyl)-thiophene-2-

carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-difluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-difluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-difluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-carbamoylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-carbamoylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-carbamoylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-sulfamoylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-sulfamoylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-sulfamoylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-

- [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-
- 5 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-difluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-difluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-difluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-
- 10 azabicyclo[2.2.2]oct-3-yl]-5-(2-carbamoylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-carbamoylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-carbamoylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-sulfamoylphenylsulfanyl)-thiophene-2-carboxamide;
- 15 N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-sulfamoylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-sulfamoylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenylsulfanyl)-
- 20 thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-
- 25 aminophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetamidophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetamidophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetamidophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-
- 30 azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoroacetamidophenoxy)-thiophene-2-

carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-difluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-difluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-difluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-carbamoylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-carbamoylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-carbamoylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-sulfamoylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-sulfamoylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-sulfamoylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-

chlorophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-morpholin-4-yl-phenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-morpholin-4-yl-phenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-morpholin-4-yl-phenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenoxy)-thiophene-2-carboxamide; N-

[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenoxy-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-morpholin-4-yl-phenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-morpholin-4-yl-phenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-morpholin-4-yl-phenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-

azabicyclo[2.2.2]oct-3-yl]-5-(3-acetylphenylsulfanyl)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetylphenylsulfanyl)-thiophene-  
2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
cyanophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
5 azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenylsulfanyl)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenylsulfanyl)-thiophene-  
2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-morpholin-4-  
yl-phenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3-morpholin-4-yl-phenylsulfanyl)-thiophene-2-  
10 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-morpholin-4-  
yl-phenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-  
methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylpyridin-4-yl)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methylpyridin-  
15 3-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-  
(5-methylpyridin-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxypyridin-4-yl)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methoxypyridin-3-yl)-thiophene-  
2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-  
20 methoxypyridin-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-yl)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-yl)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-  
2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-  
25 (5-chloropyridin-3-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-2-yl)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-2-yl)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-yloxy)-  
thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-  
30 (pyridin-3-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-2-yloxy)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylpyridin-4-yloxy)-  
thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-

methylpyridin-3-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(5-methylpyridin-2-yloxy)-thiophene-2-carboxamide; N-  
 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxypyridin-4-yloxy)-  
 thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-  
 5 methoxypyridin-3-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxypyridin-2-yloxy)-thiophene-2-carboxamide;  
 N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-yloxy)-  
 thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-  
 chloropyridin-3-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
 10 azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-yloxy)-thiophene-2-carboxamide; N-  
 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-yloxy)-  
 thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-  
 chloropyridin-3-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-yloxy)-thiophene-2-carboxamide; N-  
 15 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-3-yloxy)-  
 thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-  
 chloropyridin-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-2-yloxy)-thiophene-2-carboxamide; N-  
 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-ylsulfanyl)-thiophene-2-  
 20 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-3-  
 ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-  
 yl]-5-(pyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(2-methylpyridin-4-ylsulfanyl)-thiophene-2-carboxamide;  
 N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methylpyridin-3-ylsulfanyl)-  
 25 thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-  
 methylpyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxypyridin-4-ylsulfanyl)-thiophene-2-  
 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-  
 methoxypyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
 30 azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxypyridin-2-ylsulfanyl)-thiophene-2-  
 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-  
 4-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-  
 3-yl]-5-(6-chloropyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-

methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiophen-4-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl-thiophen-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiophen-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiophen-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(furan-4-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylfuran-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyfuran-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorofuran-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyloxazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyoxazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorooxazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-5-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methyloxazol-5-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyoxazol-5-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorooxazol-5-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiazol-2-yl)-thiophene-2-carboxamide; N-

[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-5-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylthiazol-5-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxythiazol-5-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorothiazol-5-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylthiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxythiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorothiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methyloxazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyoxazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorooxazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]thiadiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]thiadiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]thiadiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]thiadiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]oxadiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]oxadiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]oxadiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]oxadiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiophen-4-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiophen-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiophen-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiophen-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(furan-4-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylfuran-2-yloxy)-thiophene-2-carboxamide; N-

[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyfuran-2-yloxy)-  
thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-  
chlorofuran-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-  
5 2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyloxazol-2-yloxy)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-  
methoxyoxazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorooxazol-2-yloxy)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-5-yloxy)-thiophene-2-  
10 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methyloxazol-  
5-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(2-methoxyoxazol-5-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorooxazol-5-yloxy)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-2-yloxy)-thiophene-2-  
15 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiazol-  
2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(5-methoxythiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiazol-2-yloxy)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-5-yloxy)-thiophene-2-  
20 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylthiazol-  
5-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(2-methoxythiazol-5-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorothiazol-5-yloxy)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylthiazol-2-yloxy)-  
25 thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
methoxythiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorothiazol-2-yloxy)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methyloxazol-2-yloxy)-  
thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
30 methoxyoxazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorooxazol-2-yloxy)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]thiadiazol-2-yloxy)-  
thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-

methyl[1,3,4]thiadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]thiadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]thiadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]oxadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]oxadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]oxadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]oxadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiophen-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiophen-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiophen-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiophen-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(furan-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylfuran-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyfuran-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorofuran-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyloxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyoxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorooxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methyloxazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyoxazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorooxazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiazol-2-ylsulfanyl)-

thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylthiazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxythiazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorothiazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylthiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxythiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorothiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methyloxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyoxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorooxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]thiadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]thiadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]thiadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]thiadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]oxadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]oxadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]oxadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]oxadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrrole-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(isothiazol-3-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(isoxazol-3-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3H-imidazol-4-yl)-thiophene-2-

carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-(4-hydroxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-(3-acetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-(2-methanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-(pyridin-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-(2-methylpyridin-4-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-(4-trifluoromethylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-(2-acetylphenyl)-thiophene-2-carboxamide, N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-5-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-cyano-5-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-methoxy-5-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-methyl-4-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-cyano-4-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-methoxy-4-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-bromo-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-methylsulfanyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-methyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-cyano-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-bromo-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-methylsulfanyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-methyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-cyano-thiophene-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-furan-2-carboxamide; N-

[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3-dichlorophenyl)-furan-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-furan-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,5-dichlorophenyl)-furan-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,6-dichlorophenyl)-furan-2-carboxamide; N-  
5 [(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl)-furan-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-furan-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-amino-2-fluorophenyl)-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-amino-2-fluorophenyl)-furan-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-amino-2-chlorophenyl)-  
10 furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-amino-2-  
chlorophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
aminophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
cyanophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
cyanophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
15 hydroxyphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
hydroxyphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
hydroxyphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
fluoro-4-methylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-  
5-(2-fluoro-3-methoxyphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-  
20 3-yl]-5-(2-fluoro-4-methoxyphenyl)-furan-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-[(methylsulfonyl)amino]phenyl)-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-[(methylsulfonyl)amino]phenyl)-furan-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
[(methylsulfonyl)amino]phenyl)-furan-2-carboxamide; N-[(3R)-1-  
25 azabicyclo[2.2.2]oct-3-yl]-5-(2-([(trifluoromethyl)sulfonyl]amino)phenyl)-furan-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
([(trifluoromethyl)sulfonyl]amino)phenyl)-furan-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(4-([(trifluoromethyl)sulfonyl]amino)phenyl)-furan-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
30 [(phenylsulfonyl)amino]phenyl)-furan-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3-[(phenylsulfonyl)amino]phenyl)-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-[(phenylsulfonyl)amino]phenyl)-furan-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-

- [(methylamino)carbonyl]phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-[(methylamino)carbonyl]phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-[(methylamino)carbonyl]phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-[(methylamino)sulfonyl]phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-[(methylamino)sulfonyl]phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-[(methylamino)sulfonyl]phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(methylamino)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(methylamino)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(methylamino)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(ethylamino)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(ethylamino)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(ethylamino)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(methylthio)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(methylthio)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(methylthio)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(phenylthio)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(phenylthio)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(phenylthio)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-phenoxyphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-phenoxyphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-phenoxyphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-anilinophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-anilinophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-anilinophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylthio)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-fluorophenyl)thio]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-fluorophenyl)thio]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-

fluorophenyl)thio]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-chlorophenyl)thio]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-chlorophenyl)thio]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-chlorophenyl)thio]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenoxy)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenoxy)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenoxy)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenoxy)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenoxy)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-pyridin-2-yl-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-pyridin-3-yl-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-pyridin-4-yl-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyridin-2-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-2-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-2-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-2-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-3-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-3-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-3-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-3-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-4-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyridin-4-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chloropyridin-2-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-2-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-2-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-3-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-3-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-3-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chloropyridin-4-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-piperidin-1-ylphenyl)-furan-

2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-piperidin-1-ylphenyl)-  
furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-piperidin-1-  
ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
morpholin-4-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-  
5 (3-morpholin-4-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(4-morpholin-4-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-pyrrolidin-1-ylphenyl)-furan-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-pyrrolidin-1-ylphenyl)-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-pyrrolidin-1-ylphenyl)-furan-2-  
10 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(1H-pyrrol-2-yl)phenyl]-  
furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(1H-pyrrol-2-  
yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1H-  
pyrrol-2-yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-  
(2-furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(3-  
15 furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(5-  
methyl-2-furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-  
5-[3-(2-furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-  
[3-(3-furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-  
(5-methyl-2-furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
20 yl]-5-[4-(2-furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-  
5-[4-(3-furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-  
[4-(5-methyl-2-furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-  
3-yl]-5-[2-(1,3-thiazol-2-yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-[3-(1,3-thiazol-2-yl)phenyl]-furan-2-carboxamide; N-  
25 [(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1,3-thiazol-2-yl)phenyl]-furan-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(1,3-oxazol-2-yl)phenyl]-  
furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(1,3-oxazol-2-  
yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1,3-  
oxazol-2-yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
30 isothiazol-5-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-  
(3-isothiazol-5-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-  
5-(4-isothiazol-5-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-[4-(1H-indol-2-yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-

- azabicyclo[2.2.2]oct-3-yl]-5-[4-(1H-indol-3-yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(1H-indol-5-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(1H-indol-6-yl)-furan-2-carboxamide;
- 5 N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3-dichlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-
- 10 azabicyclo[2.2.2]oct-3-yl]-5-(2,5-dichlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,6-dichlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-
- 15 3-yl]-5-(4-amino-2-fluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-amino-2-fluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-amino-2-chlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-amino-2-chlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-
- 20 3-yl]-5-(3-aminophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-
- 25 hydroxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoro-4-methylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoro-3-methoxyphenyl)-furan-
- 30 methoxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-[(methylsulfonyl)amino]phenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-[(methylsulfonyl)amino]phenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-

[(methylsulfonyl)amino]phenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-[2-([(trifluoromethyl)sulfonyl]amino)phenyl]-furan-2-  
 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-  
 ([trifluoromethyl)sulfonyl]amino)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-  
 5 methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-([(trifluoromethyl)sulfonyl]amino)phenyl)-  
 furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
 [(phenylsulfonyl)amino]phenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(3-[(phenylsulfonyl)amino]phenyl)-furan-2-carboxamide;  
 N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
 10 [(phenylsulfonyl)amino]phenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(2-[(methylamino)carbonyl]phenyl)-furan-2-  
 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
 [(methylamino)carbonyl]phenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(4-[(methylamino)carbonyl]phenyl)-furan-2-  
 15 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
 [(methylamino)sulfonyl]phenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(3-[(methylamino)sulfonyl]phenyl)-furan-2-carboxamide;  
 N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
 [(methylamino)sulfonyl]phenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
 20 azabicyclo[2.2.2]oct-3-yl]-5-[2-(methylamino)phenyl]-furan-2-carboxamide; N-  
 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(methylamino)phenyl]-furan-2-  
 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-  
 (methylamino)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-[2-(ethylamino)phenyl]-furan-2-carboxamide; N-  
 25 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(ethylamino)phenyl]-furan-2-  
 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-  
 (ethylamino)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(2-acetylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-  
 methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetylphenyl)-furan-2-carboxamide; N-  
 30 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetylphenyl)-furan-2-  
 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-  
 (methylthio)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-[3-(methylthio)phenyl]-furan-2-carboxamide; N-

[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(methylthio)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(phenylthio)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(phenylthio)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(phenylthio)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-phenoxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-phenoxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-phenoxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-anilinophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-anilinophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-anilinophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylthio)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-fluorophenyl)thio]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-fluorophenyl)thio]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-fluorophenyl)thio]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-chlorophenyl)thio]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-chlorophenyl)thio]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-chlorophenyl)thio]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenoxy)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenoxy)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenoxy)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenoxy)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenoxy)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-pyridin-2-yl-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-pyridin-3-yl-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-pyridin-4-yl-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyridin-2-yl)-furan-2-carboxamide;

N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-2-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-2-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-2-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-3-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-3-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-3-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-3-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-4-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyridin-4-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chloropyridin-2-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-2-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-2-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-3-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-3-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-3-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chloropyridin-4-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-piperidin-1-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-piperidin-1-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-piperidin-1-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-morpholin-4-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-morpholin-4-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-morpholin-4-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-pyrrolidin-1-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-

(3-pyrrolidin-1-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-pyrrolidin-1-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(1H-pyrrol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(1H-pyrrol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1H-pyrrol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(2-furyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(3-furyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(5-methyl-2-furyl)phenyl]-furan-2-carboxamid; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(2-furyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(3-furyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(5-methyl-2-furyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(2-furyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(3-furyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(5-methyl-2-furyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(1,3-thiazol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(1,3-thiazol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1,3-thiazol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(1,3-oxazol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(1,3-oxazol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1,3-oxazol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-isothiazol-5-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-isothiazol-5-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-isothiazol-5-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1H-indol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1H-indol-3-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(1H-indol-5-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(1H-indol-6-yl)-furan-2-carboxamide;

- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2,3-difluorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2,4-difluorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2,5-difluorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2,6-difluorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3,4-difluorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3,5-difluorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-chlorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-chlorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-chlorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-bromophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-bromophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-bromophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-cyanophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-cyanophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-cyanophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-nitrophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-nitrophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-nitrophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-methylphenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-methylphenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-aminophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-aminophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-aminophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[2-(methylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[3-(methylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[4-(methylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[2-(acetylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[3-(acetylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[4-(acetylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-

[(trifluoroacetyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-[(trifluoroacetyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-[(trifluoroacetyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-[(methylsulfonyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-[(methylsulfonyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-[(methylsulfonyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-hydroxyphenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-hydroxyphenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-methoxyphenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-methoxyphenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-methoxyphenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[2-(trifluoromethoxy)phenyl]-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[3-(trifluoromethoxy)phenyl]-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[4-(trifluoromethoxy)phenyl]-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-pyridin-2-yl-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-pyridin-3-yl-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-pyridin-4-yl-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(6-fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(5-fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(6-fluoropyridin-3-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(5-fluoropyridin-3-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-fluoropyridin-3-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-fluoropyridin-3-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-fluoropyridin-4-

yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-fluoropyridin-4-yl)-1,3-thiazole-5-carboxamide;

- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(methylthio)-1,3-thiazole-2-carboxamide; N-  
5 [(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylthio)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-fluorophenyl)thio]-1,3-thiazole-2-carboxamide; N-  
10 [(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-fluorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-fluorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-chlorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-chlorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-chlorophenyl)thio]-1,3-thiazole-2-carboxamide; N-  
15 [(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenoxy)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenoxy)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenoxy)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenoxy)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenoxy)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-

5-(2-hydroxyphenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-(acetylamino)phenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(acetylamino)phenyl]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetylamino)phenyl]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-4-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-methoxyphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-methoxyphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-methoxyphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-methylphenyl)thio]-1,3-thiazole-2-carboxamide; N-

[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-methylphenyl)thio]-1,3-thiazole-2-  
 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-methylphenyl)thio]-1,3-  
 thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-([4-  
 (acetylamino)phenyl]thio)-1,3-thiazole-2-carboxamide; N-[(3R)-1-  
 5 azabicyclo[2.2.2]oct-3-yl]-5-[(4-aminophenyl)thio]-1,3-thiazole-2-carboxamide; N-  
 [(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-hydroxyphenyl)thio]-1,3-thiazole-2-  
 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenoxy)-1,3-  
 thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenoxy)-  
 1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
 10 methylphenoxy)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-  
 (2-methoxyphenoxy)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
 yl]-5-(3-methoxyphenoxy)-1,3-thiazole-2-carboxamide; N-[(3R)-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenoxy)-1,3-thiazole-2-carboxamide; N-  
 [(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenoxy)-1,3-thiazole-2-  
 15 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenoxy)-1,3-  
 thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-  
 (acetylamino)phenoxy]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-  
 3-yl]-5-(2-fluorophenyl)-4-methyl-1,3-thiazole-2-carboxamide; N-[(3R)-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-4-methyl-1,3-thiazole-2-carboxamide;  
 20 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-4-methyl-1,3-thiazole-2-  
 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-4-methyl-1,3-  
 thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-  
 4-methyl-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
 methoxyphenyl)-4-methyl-1,3-thiazole-2-carboxamide; N-[(3R)-1-  
 25 azabicyclo[2.2.2]oct-3-yl]-4-methyl-5-thien-2-yl-1,3-thiazole-2-carboxamide,  
  
 N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2,3-difluorophenyl)-1,3-  
 thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2,4-  
 difluorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-  
 30 azabicyclo[2.2.2]oct-3-yl]-2-(2,5-difluorophenyl)-1,3-thiazole-5-carboxamide; N-  
 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2,6-difluorophenyl)-1,3-thiazole-5-  
 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3,4-  
 difluorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-

azabicyclo[2.2.2]oct-3-yl]-2-(3,5-difluorophenyl)-1,3-thiazole-5-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-chlorophenyl)-1,3-thiazole-5-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-chlorophenyl)-  
1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-  
5 chlorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-2-(2-bromophenyl)-1,3-thiazole-5-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-bromophenyl)-1,3-thiazole-5-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-bromophenyl)-  
1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-  
10 cyanophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-2-(3-cyanophenyl)-1,3-thiazole-5-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-cyanophenyl)-1,3-thiazole-5-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-nitrophenyl)-  
1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-  
15 nitrophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-2-(4-nitrophenyl)-1,3-thiazole-5-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-methylphenyl)-1,3-thiazole-5-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-methylphenyl)-  
1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-  
20 aminophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-2-(3-aminophenyl)-1,3-thiazole-5-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-aminophenyl)-1,3-thiazole-5-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[2-  
(methylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-  
25 azabicyclo[2.2.2]oct-3-yl]-2-[3-(methylamino)phenyl]-1,3-thiazole-5-carboxamide;  
N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[4-(methylamino)phenyl]-1,3-  
thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[2-  
(acetylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-2-[3-(acetylamino)phenyl]-1,3-thiazole-5-carboxamide; N-  
30 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[4-(acetylamino)phenyl]-1,3-  
thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-  
[(trifluoroacetyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-2-(3-[(trifluoroacetyl)amino]phenyl)-1,3-thiazole-5-

carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-  
[(trifluoroacetyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-2-(2-[(methylsulfonyl)amino]phenyl)-1,3-thiazole-5-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-  
5 [(methylsulfonyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-  
1-azabicyclo[2.2.2]oct-3-yl]-2-(4-[(methylsulfonyl)amino]phenyl)-1,3-thiazole-5-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-  
hydroxyphenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-2-(4-hydroxyphenyl)-1,3-thiazole-5-carboxamide; N-  
10 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-methoxyphenyl)-1,3-thiazole-5-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-  
methoxyphenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-2-(4-methoxyphenyl)-1,3-thiazole-5-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[2-(trifluoromethoxy)phenyl]-1,3-  
15 thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[3-  
(trifluoromethoxy)phenyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-2-[4-(trifluoromethoxy)phenyl]-1,3-thiazole-5-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-pyridin-2-yl-1,3-  
thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-pyridin-  
20 3-yl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-  
2-pyridin-4-yl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-2-(6-fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(5-fluoropyridin-2-yl)-1,3-thiazole-  
5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-  
25 fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-2-(3-fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-fluoropyridin-2-yl)-1,3-thiazole-  
5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(6-  
fluoropyridin-3-yl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-  
30 azabicyclo[2.2.2]oct-3-yl]-2-(5-fluoropyridin-3-yl)-1,3-thiazole-5-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-fluoropyridin-3-yl)-1,3-thiazole-  
5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-  
fluoropyridin-3-yl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-

azabicyclo[2.2.2]oct-3-yl]-2-(2-fluoropyridin-4-yl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-fluoropyridin-4-yl)-1,3-thiazole-5-carboxamide;

- 5 N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(methylthio)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-1,3-thiazole-2-carboxamide, N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylthio)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-fluorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-fluorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-fluorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-chlorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-chlorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-chlorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-thien-3-yl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-

1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(acetylamino)phenyl]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(acetylamino)phenyl]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetylamino)phenyl]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-

fluoropyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-4-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-methoxyphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-methoxyphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-methoxyphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-methylphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-methylphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-methylphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-(acetylamino)phenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-aminophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-hydroxyphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetylamino)phenoxy]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-4-methyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-4-methyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-

azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-4-methyl-1,3-thiazole-2-carboxamide;  
N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-4-methyl-1,3-  
thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
methoxyphenyl)-4-methyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
5 azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-4-methyl-1,3-thiazole-2-  
carboxamide;  
N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-5-thien-2-yl-1,3-thiazole-  
2-carboxamide, N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1,3-  
thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-  
10 4-methyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-  
3-yl]-5-(2-fluorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1,3-thiazole-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1,3-thiazole-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-thien-2-yl-1,3-  
15 thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
chlorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1,3-thiazole-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-1,3-thiazole-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-  
20 1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
nitrophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1,3-thiazole-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-1,3-thiazole-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
25 methoxyphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-2-(2-hydroxyphenyl)-1,3-thiazole-5-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-methylphenyl)-1,3-thiazole-5-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-phenyl-1,3-  
thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-  
30 fluorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-2-(3-fluorophenyl)-1,3-thiazole-5-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-fluorophenyl)-1,3-thiazole-5-  
carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenylsulfanyl-1,3-oxazole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenylsulfanyl)-1,3-oxazole-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenylsulfanyl)-1,3-  
5 oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
fluorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(3,4-difluorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenylsulfanyl)-1,3-oxazole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenylsulfanyl)-1,3-oxazole-2-  
10 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenylsulfanyl)-1,3-  
oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
chlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(4-chlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenylsulfanyl)-1,3-oxazole-2-  
15 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenylsulfanyl)-  
1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-  
dichlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-  
3-yl]-5-(2-fluorophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-  
20 1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenyl)-1,3-oxazole-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-1,3-oxazole-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-1,3-  
oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-  
1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-  
25 dichlorophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-  
(2,4-dichlorophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(3,5-dichlorophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-  
1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-1,3-oxazole-2-carboxamide; N-  
30 [(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-1,3-oxazole-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-1,3-oxazole-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-1,3-oxazole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-1,3-oxazole-2-

carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylaminophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylaminophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylaminophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(acetylamino)phenyl]-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(acetylamino)phenyl]-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetylamino)phenyl]-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-2-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-3-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyridin-2-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-2-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-2-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-2-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-3-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-3-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-3-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-3-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-4-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-3-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-

5-(3-fluorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenoxy)-1,3-oxazole-2-carboxamide;

N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenylsulfanyl-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1,3-oxazole-2-

carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylaminophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylaminophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylaminophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(acetylamino)phenyl]-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-

azabicyclo[2.2.2]oct-3-yl]-5-[3-(acetylamino)phenyl]-1,3-oxazole-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetylamino)phenyl]-1,3-  
oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-  
(pyridin-2-yl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
5 azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-3-yl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-  
2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-yl)-1,3-oxazole-2-carboxamide;  
N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyridin-2-yl)-1,3-  
oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
fluoropyridin-2-yl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
10 azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-2-yl)-1,3-oxazole-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-2-yl)-1,3-oxazole-  
2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
fluoropyridin-3-yl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-3-yl)-1,3-oxazole-2-carboxamide; N-  
15 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-3-yl)-1,3-oxazole-  
2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-  
fluoropyridin-3-yl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-4-yl)-1,3-oxazole-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-3-yl)-1,3-oxazole-  
20 2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-1,3-  
oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
fluorophenoxy)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenoxy)-1,3-oxazole-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenoxy)-1,3-oxazole-2-  
25 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-  
difluorophenoxy)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenoxy)-1,3-oxazole-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenoxy)-1,3-oxazole-  
2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
30 chlorophenoxy)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenoxy)-1,3-oxazole-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-1,3-oxazole-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-

dichlorophenoxy)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenoxy)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenoxy)-1,3-oxazole-2-carboxamide;

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N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(methylthio)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(methylthio)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,6-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4,6-trifluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-

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azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylaminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylaminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylaminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(acetyl amino)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(acetyl amino)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetyl amino)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-2-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyrid-2-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyrid-2-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyrid-2-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyrid-2-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-

azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyrid-4-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyrid-4-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-1H-pyrrole-2-carboxamide; N-

[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-1H-pyrrole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-1H-pyrrole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-1H-pyrrole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylaminophenyl)-1H-pyrrole-2-  
5 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylaminophenyl)-1H-  
pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
methylaminophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(2-nitrophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(3-nitrophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
10 yl]-5-(4-nitrophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(2-aminophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(3-aminophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(4-aminophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-[2-(acetylamino)phenyl]-1H-pyrrole-2-carboxamide; N-[(3R)-1-  
15 azabicyclo[2.2.2]oct-3-yl]-5-[3-(acetylamino)phenyl]-1H-pyrrole-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetylamino)phenyl]-1H-pyrrole-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-2-yl)-1H-pyrrole-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-3-yl)-1H-pyrrole-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-4-yl)-1H-pyrrole-2-  
20 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyrid-2-yl)-1H-  
pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyrid-2-yl)-  
1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyrid-2-  
yl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-  
fluoropyrid-2-yl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-  
25 (2-fluoropyrid-3-yl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(4-fluoropyrid-3-yl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyrid-3-yl)-1H-pyrrole-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyrid-3-yl)-1H-pyrrole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyrid-4-yl)-1H-pyrrole-2-  
30 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyrid-4-yl)-1H-  
pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-1H-  
pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-1H-

pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-1H-pyrrole-2-carboxamide;

- 5 N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(methylthio)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(methylthio)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,6-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4,6-trifluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-1-methyl-1H-pyrrole-2-carboxamide;

N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylaminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylaminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylaminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(acetyl amino)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(acetyl amino)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetyl amino)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-2-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-pyridin-4-yl-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyrid-2-yl)-1-methyl-1H-

pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyrid-2-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyrid-2-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyrid-2-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyrid-4-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyrid-4-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-1H-pyrrole-2-carboxamide; N-

[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylaminophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylaminophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylaminophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(acetylamino)phenyl]-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(acetylamino)phenyl]-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetylamino)phenyl]-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-2-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-3-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-

azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-4-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyrid-2-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyrid-2-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyrid-2-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyrid-2-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyrid-3-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyrid-3-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyrid-3-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyrid-3-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyrid-4-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyrid-4-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-1H-pyrrole-2-carboxamide; or N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-1H-pyrrole-2-carboxamide.

## Materials and Methods for identifying binding constants:

**Membrane Preparation.** Male Sprague-Dawley rats (300-350g) are sacrificed by decapitation and the brains (whole brain minus cerebellum) are dissected quickly, weighed and homogenized in 9 volumes/g wet weight of ice-cold 0.32 M sucrose using a rotating pestle on setting 50 (10 up and down strokes). The homogenate is centrifuged at 1,000 x g for 10 minutes at 4 °C. The supernatant is collected and centrifuged at 20,000 x g for 20 minutes at 4 °C. The resulting pellet is resuspended to a protein concentration of 1 - 8 mg/mL. Aliquots of 5 mL homogenate are frozen at -80 °C until needed for the assay. On the day of the assay, aliquots are thawed at room temperature and diluted with Krebs' - 20 mM Hepes buffer pH 7.0 (at room temperature) containing 4.16 mM NaHCO<sub>3</sub>, 0.44 mM KH<sub>2</sub>PO<sub>4</sub>, 127 mM NaCl, 5.36 mM KCl, 1.26 mM CaCl<sub>2</sub>, and 0.98 mM MgCl<sub>2</sub>, so that 25 - 150 µg protein are added per test tube. Proteins are determined by the Bradford method (Bradford, M.M., *Anal. Biochem.*, 72, 248-254, 1976) using bovine serum albumin as the standard.

Binding Assay. For saturation studies, 0.4 mL homogenate are added to test tubes containing buffer and various concentrations of radioligand, and are incubated in a final volume of 0.5 mL for 1 hour at 25 °C. Nonspecific binding was determined in tissues incubated in parallel in the presence of 0.05 ml MLA for a final  
5 concentration of 1  $\mu$ M MLA, added before the radioligand. In competition studies, drugs are added in increasing concentrations to the test tubes before addition of 0.05 ml [ $^3$ H]-MLA for a final concentration of 3.0 to 4.0 nM [ $^3$ H]-MLA. The incubations are terminated by rapid vacuum filtration through Whatman GF/B glass filter paper mounted on a 48 well Brandel cell harvester. Filters are pre-soaked in 50 mM Tris  
10 HCl pH 7.0 - 0.05 % polyethylenimine. The filters are rapidly washed two times with 5 mL aliquots of cold 0.9% saline and then counted for radioactivity by liquid scintillation spectrometry.

Data Analysis. In competition binding studies, the inhibition constant ( $K_i$ ) was calculated from the concentration dependent inhibition of [ $^3$ H]-MLA binding  
15 obtained from non-linear regression fitting program according to the Cheng-Prusoff equation (Cheng, Y.C. and Prusoff, W.H., *Biochem. Pharmacol.*, 22, p. 3099-3108, 1973). Hill coefficients were obtained using non-linear regression (GraphPad Prism sigmoidal dose-response with variable slope).

The aforementioned examples have the provided Ki values:

Example #	Ki (nM)
1	13-33
2	110
3	95-157
4	60-73
5	9-15
6	5-10
7	25-30
8	46
9	540-570
12	45-75
14	12-16
15	40-90
16	3-8
17	260-385
18	240-290
19	2-4
20	3-5
21	1-2
22	1-6
23	6-10
25	10-15
26	5-20
27	35-160
33	27
34	30-39
35	30-50
36	110-115
40	12-16
41	3-4
43	3-5

Example #	Ki (nM)
50	68-564
52	340-480
53	1900-1970
55	60-70
58	920-1005
79	20-75
80	50-60
81	55-175
82	115-125
83	2-3
84	1100-1490
85	30-40
86	30-40
87	130-150
88	65-80
89	310-400
90	80-150
92	190-375
93	40-45
94	125
95	1085
96	95-110
98	50-80
99	310-325
103	17-50
104	20-130
104	75
105	90
106	15-190
106	22-110

30

Example #	Ki (nM)
108	415-560
109	1.5-6
110	4-190
112	5-24
113	90-125
114	135
117	1860
118	45-69
119	415-560
120	1050
128	6-7
130	420-565
130	330-410
132	875-2085
133	30-70
134	60-100
135	16-45
136	15-20
138	525-590
140	35-65
141	260-275
146	35-140
147	2-3
150	290-330
151	225-835
152	20-65
153	1555
154	>1000
155	1380
158	12-20

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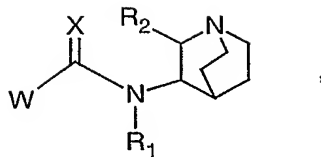
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## CLAIMS

What is claimed:

1. A compound of formula I:



Formula I

or pharmaceutically acceptable salts thereof, wherein

R<sub>1</sub> is selected from -H, alkyl, cycloalkyl, halogenated alkyl, or aryl;

Alkyl is both straight and branched-chain moieties having from 1-6 carbon atoms;

Halogenated alkyl is an alkyl moiety having from 1-6 carbon atoms and having 1 to (2n+1) substituent(s) independently selected from -F, -Cl, -Br, or -I where n is the maximum number of carbon atoms in the moiety;

Cycloalkyl is a cyclic alkyl moiety having from 3-6 carbon atoms;

Aryl is phenyl, substituted phenyl, naphthyl, or substituted naphthyl;

Substituted phenyl is a phenyl either having 1-4 substituents independently selected from -F, -Cl, -Br, or -I, or having 1 substituent selected from -R<sub>12</sub> and 0-3 substituents independently selected from -F, -Cl, -Br, or -I;

Substituted naphthyl is a naphthalene moiety either having 1-4 substituents independently selected from -F, -Cl, -Br, or -I, or having 1 substituent selected from -R<sub>12</sub> and 0-3 substituents independently selected from -F, -Cl, -Br, or -I, where the substitution can be independently on either the same ring or different rings of said naphthalene moiety;

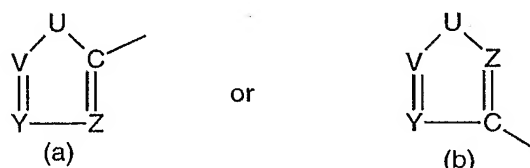
R<sub>2</sub> is -H, alkyl, halogenated alkyl, substituted alkyl, cycloalkyl, benzyl, substituted benzyl, or aryl;

Substituted alkyl is an alkyl moiety having from 1-6 carbon atoms and having 0-3 substituents independently selected from -F, -Cl, -Br, or -I, and further having 1 substituent selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -NO<sub>2</sub>, -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl;

Substituted benzyl is a benzyl either having 1-4 substituents independently selected from -F, -Cl, -Br, or -I, or having 1 substituent selected from -R<sub>12</sub> and 0-3 substituents independently selected from -F, -Cl, -Br, or -I;

X is O or S;

- 5 W is a cyclic heteroaromatic moiety where the heteroatoms can be from 1-3 atoms selected from oxygen, sulfur, or nitrogen of the following structures:



wherein U is -O-, -S-, or -N(R<sub>3</sub>)-;

V and Y are independently selected from =N-, or =C(R<sub>5</sub>)-;

- 10 Z is =N-, or =CH-, provided that when both V and Y are =C(R<sub>5</sub>)- and Z is =CH-, only one =C(R<sub>5</sub>)- can be =CH-, and further provided that when U is -O-, Y is =C(R<sub>5</sub>)- and Z is =C(R<sub>5</sub>)-, V cannot be =N-;

- R<sub>3</sub> is -H, alkyl, cycloalkyl, heterocycloalkyl, halogenated alkyl, halogenated cycloalkyl, halogenated heterocycloalkyl, substituted alkyl, substituted cycloalkyl, substituted heterocycloalkyl, or aryl, and provided that when W is (b) and Z is =N- and U is N(R<sub>3</sub>), R<sub>3</sub> cannot be phenyl or substituted phenyl
- 15

Alkenyl is straight and branched-chain moieties having from 2-6 carbon atoms and having at least one carbon-carbon double bond;

- Halogenated alkenyl is an unsaturated alkenyl moiety having from 2-6 carbon atoms and having 1 to (2n-1) substituent(s) independently selected from -F, -Cl, -Br, or -I where n is the maximum number of carbon atoms in the moiety;
- 20

- Substituted alkenyl is an unsaturated alkenyl moiety having from 2-6 carbon atoms and having 0-3 substituents independently selected from -F, or -Cl, and further having 1 substituent selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl;
- 25

Alkynyl is straight and branched-chain moieties having from 2-6 carbon atoms and having at least one carbon-carbon triple bond;

Halogenated alkynyl is an unsaturated alkynyl moiety having from 3-6 carbon atoms and having 1 to (2n-3) substituent(s) independently selected from -F, -Cl, -Br, or -I where n is the maximum number of carbon atoms in the moiety;

Substituted alkynyl is an unsaturated alkynyl moiety having from 3-6 carbon atoms and having 0-3 substituents independently selected from -F, or -Cl, and further having 1 substituent selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -CN, -C(O)NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl;

Halogenated cycloalkyl is a cyclic moiety having from 3-6 carbon atoms and having 1-4 substituents independently selected from -F, or -Cl;

Substituted cycloalkyl is a cyclic moiety having from 3-6 carbon atoms and having 0-3 substituents independently selected from -F, or -Cl, and further having 1 substituent selected from -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, -NO<sub>2</sub>, phenyl, or substituted phenyl;

Heterocycloalkyl is a cyclic moiety having 4-7 atoms with 1-2 atoms within the ring being -S-, -N(R<sub>3</sub>)-, or -O-;

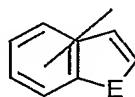
Halogenated heterocycloalkyl is a cyclic moiety having from 4-7 atoms with 1-2 atoms within the ring being -S-, -N(R<sub>3</sub>)-, or -O-, and having 1-4 substituents independently selected from -F, or -Cl;

Substituted heterocycloalkyl is a cyclic moiety having from 4-7 atoms with 1-2 atoms within the ring being -S-, -N(R<sub>3</sub>)-, or -O- and having 0-3 substituents independently selected from -F, or -Cl, and further having 1 substituent selected from -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -NO<sub>2</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl;

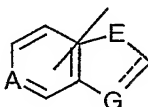
R<sub>5</sub> is independently selected from the group consisting of -H, alkyl, alkenyl, alkynyl, cycloalkyl, heterocycloalkyl, halogenated alkyl, halogenated alkenyl, halogenated alkynyl, halogenated cycloalkyl, halogenated heterocycloalkyl, substituted alkyl, substituted alkenyl, substituted alkynyl, substituted cycloalkyl, substituted heterocycloalkyl, aryl, -OR<sub>8</sub>, -SR<sub>8</sub>, -F, -Cl, -Br, -I, -NR<sub>8</sub>R<sub>8</sub>, -C(O)R<sub>8</sub>, -C(O)NR<sub>8</sub>R<sub>8</sub>, -CN, -NR<sub>8</sub>C(O)R<sub>11</sub>, -S(O)<sub>2</sub>NR<sub>8</sub>R<sub>8</sub>, -OS(O)<sub>2</sub>R<sub>11</sub>, -S(O)<sub>2</sub>R<sub>8</sub>, -NR<sub>8</sub>S(O)<sub>2</sub>R<sub>8</sub>, -N(H)C(O)N(H)R<sub>8</sub>, -NO<sub>2</sub>, -R<sub>7</sub>, and -R<sub>9</sub>;

R<sub>7</sub> is 5-membered heteroaromatic mono-cyclic moieties containing within the ring 1-3 heteroatoms independently selected from the group consisting of -O-, =N-,

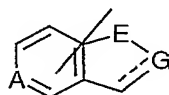
-N(R<sub>3</sub>)-, and -S-, and having 0-1 substituent selected from -R<sub>12</sub> and 0-3 substituents independently selected from -F, -Cl, -Br, or -I, or R<sub>7</sub> is a 9-membered fused-ring moiety having a 6-membered ring fused to a 5-membered ring and having the formula



5 wherein E is O, S, or NR<sub>3</sub>,



wherein E and G are independently selected from CR<sub>18</sub>, O, S, or NR<sub>3</sub>, and A is CR<sub>18</sub> or N, or



10 wherein E and G are independently selected from CR<sub>18</sub>, O, S, or NR<sub>3</sub>, and A is CR<sub>18</sub> or N, each 9-membered fused-ring moiety having 0-1 substituent selected from -R<sub>12</sub> and 0-3 substituent(s) independently selected from -F, -Cl, -Br, or -I, and having a bond directly or indirectly attached to the core molecule where valency allows in either the 6-membered or the 5-membered ring of the fused-ring moiety;

15 Each R<sub>8</sub> is independently selected from -H, alkyl, halogenated alkyl, substituted alkyl, cycloalkyl, halogenated cycloalkyl, substituted cycloalkyl, heterocycloalkyl, halogenated heterocycloalkyl, substituted heterocycloalkyl, -R<sub>7</sub>, -R<sub>9</sub>, phenyl, or substituted phenyl;

R<sub>9</sub> is 6-membered heteroaromatic mono-cyclic moieties containing within the  
 20 ring 1-3 heteroatoms selected from =N- and having 0-1 substituent selected from -R<sub>12</sub> and 0-3 substituent(s) independently selected from -F, -Cl, -Br, or -I, or 10-membered heteroaromatic bi-cyclic moieties containing within one or both rings 1-3 heteroatoms selected from =N-, including quinolinyl or isoquinolinyl, each 10-membered fused-ring moiety having 0-1 substituent selected from -R<sub>12</sub> and 0-3 substituent(s)  
 25 independently selected from -F, -Cl, -Br, or -I and having a bond directly or indirectly attached to the core molecule where valency allows;

Each R<sub>10</sub> is independently selected from -H, alkyl, cycloalkyl, heterocycloalkyl, alkyl substituted with 1 substituent selected from R<sub>13</sub>, cycloalkyl substituted with 1 substituent selected from R<sub>13</sub>, heterocycloalkyl substituted with 1

substituent selected from  $R_{13}$ , halogenated alkyl, halogenated cycloalkyl, halogenated heterocycloalkyl, phenyl, substituted phenyl,  $-R_7$ , or  $-R_9$ ;

Each  $R_{11}$  is independently selected from  $-H$ , alkyl, cycloalkyl, heterocycloalkyl, halogenated alkyl, halogenated cycloalkyl, or halogenated heterocycloalkyl;

5  $R_{12}$  is selected from  $-OR_{11}$ ,  $-SR_{11}$ , alkyl, cycloalkyl, heterocycloalkyl, halogenated alkyl, halogenated cycloalkyl, halogenated heterocycloalkyl, substituted alkyl, substituted cycloalkyl, substituted heterocycloalkyl,  $-NR_{11}R_{11}$ ,  $-C(O)R_{11}$ ,  $-NO_2$ ,  $-C(O)NR_{11}R_{11}$ ,  $-CN$ ,  $-NR_{11}C(O)R_{11}$ ,  $-S(O)_2NR_{11}R_{11}$ , or  $-NR_{11}S(O)_2R_{11}$ ;

10  $R_{13}$  is selected from  $-OR_{11}$ ,  $-SR_{11}$ ,  $-NR_{11}R_{11}$ ,  $-C(O)R_{11}$ ,  $-C(O)NR_{11}R_{11}$ ,  $-CN$ ,  $-NR_{11}C(O)R_{11}$ ,  $-S(O)_2NR_{11}R_{11}$ ,  $-NR_{11}S(O)_2R_{11}$ ,  $-CF_3$ , or  $-NO_2$ ;

$R_{14}$  is independently selected from  $-H$ , alkyl, halogenated alkyl, limited substituted alkyl, cycloalkyl, halogenated cycloalkyl, substituted cycloalkyl, heterocycloalkyl, halogenated heterocycloalkyl, substituted heterocycloalkyl;

Each  $R_{18}$  is independently selected from  $-H$ , alkyl, cycloalkyl,  
 15 heterocycloalkyl, halogenated alkyl, halogenated cycloalkyl, halogenated heterocycloalkyl, substituted alkyl, substituted cycloalkyl, substituted heterocycloalkyl,  $-OR_{11}$ ,  $-SR_{11}$ ,  $-NR_{11}R_{11}$ ,  $-C(O)R_{11}$ ,  $-NO_2$ ,  $-C(O)NR_{11}R_{11}$ ,  $-CN$ ,  $-NR_{11}C(O)R_{11}$ ,  $-S(O)_2NR_{11}R_{11}$ , or  $-NR_{11}S(O)_2R_{11}$ ,  $-F$ ,  $-Cl$ ,  $-Br$ , or  $-I$ , or a bond directly or indirectly attached to the core molecule, provided that there is only one said  
 20 bond to the core molecule within the 9-membered fused-ring moiety, further provided that the fused-ring moiety has 0-1 substituent selected from alkyl, cycloalkyl, heterocycloalkyl, halogenated alkyl, halogenated cycloalkyl, halogenated heterocycloalkyl, substituted alkyl, substituted cycloalkyl, substituted heterocycloalkyl,  $-OR_{11}$ ,  $-SR_{11}$ ,  $-NR_{11}R_{11}$ ,  $-C(O)R_{11}$ ,  $-NO_2$ ,  $-C(O)NR_{11}R_{11}$ ,  $-CN$ ,  
 25  $-NR_{11}C(O)R_{11}$ ,  $-S(O)_2NR_{11}R_{11}$ , or  $-NR_{11}S(O)_2R_{11}$ , and further provided that the fused-ring moiety has 0-3 substituent(s) selected from  $-F$ ,  $-Cl$ ,  $-Br$ , or  $-I$ .

Limited substituted alkyl is a substituted alkyl having from 1-6 carbon atoms and having 0-3 substituents independently selected from  $-F$ ,  $-Cl$ ,  $-Br$ , or  $-I$ , and further having 1 substituent on either only the  $\omega$  carbon and selected from  $-OR_{11}$ ,  $-SR_{11}$ ,  
 30  $-NR_{11}R_{11}$ ,  $-C(O)R_{11}$ ,  $-NO_2$ ,  $-C(O)NR_{11}R_{11}$ ,  $-CN$ ,  $-NR_{10}C(O)R_{11}$ ,  $-S(O)_2NR_{10}R_{10}$ , or  $-NR_{10}S(O)_2R_{10}$ , or on any carbon with sufficient valency but not on the  $\omega$  carbon and selected from  $-R_7$ ,  $-R_9$ ,  $-OR_{10}$ ,  $-SR_{10}$ ,  $-NR_{10}R_{10}$ ,  $-C(O)R_{10}$ ,  $-NO_2$ ,  $-C(O)NR_{10}R_{10}$ ,  $-CN$ ,  $-NR_{10}C(O)R_{10}$ ,  $-S(O)_2NR_{10}R_{10}$ ,  $-NR_{10}S(O)_2R_{10}$ , phenyl, or substituted phenyl;

Limited substituted alkenyl is a substituted alkenyl having from 1-6 carbon atoms and having 0-3 substituents independently selected from -F, -Cl, -Br, or -I, and further having 1 substituent on either only the  $\omega$  carbon and selected from -OR<sub>11</sub>, -SR<sub>11</sub>, -NR<sub>11</sub>R<sub>11</sub>, -C(O)R<sub>11</sub>, -NO<sub>2</sub>, -C(O)NR<sub>11</sub>R<sub>11</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>11</sub>,  
 5 -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, or -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, or on any carbon with sufficient valency but not on the  $\omega$  carbon and selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -NO<sub>2</sub>, -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl; and

Limited substituted alkynyl is a substituted alkynyl having from 1-6 carbon  
 10 atoms and having 0-3 substituents independently selected from -F, -Cl, -Br, or -I, and further having 1 substituent on either only the  $\omega$  carbon and selected from -OR<sub>11</sub>, -SR<sub>11</sub>, -NR<sub>11</sub>R<sub>11</sub>, -C(O)R<sub>11</sub>, -NO<sub>2</sub>, -C(O)NR<sub>11</sub>R<sub>11</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>11</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, or -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, or on any carbon with sufficient valency but not on the  $\omega$  carbon and selected from -R<sub>7</sub>, -R<sub>9</sub>, -OR<sub>10</sub>, -SR<sub>10</sub>, -NR<sub>10</sub>R<sub>10</sub>, -C(O)R<sub>10</sub>, -NO<sub>2</sub>,  
 15 -C(O)NR<sub>10</sub>R<sub>10</sub>, -CN, -NR<sub>10</sub>C(O)R<sub>10</sub>, -S(O)<sub>2</sub>NR<sub>10</sub>R<sub>10</sub>, -NR<sub>10</sub>S(O)<sub>2</sub>R<sub>10</sub>, phenyl, or substituted phenyl.

2. The compounds according to claim 1, wherein X is O.

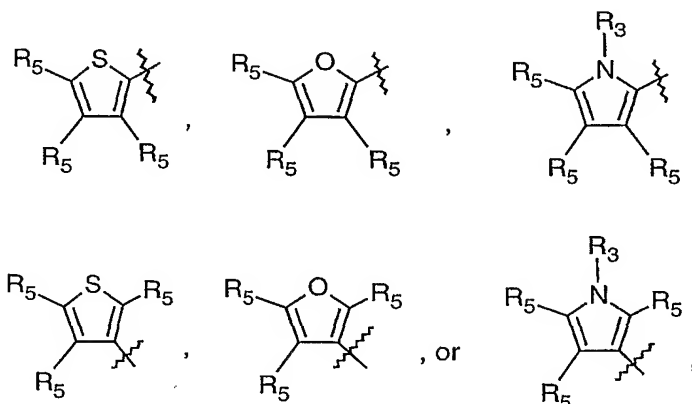
3. The compounds according to claim 2, wherein R<sub>2</sub> is one of alkyl, halogenated  
 20 alkyl, substituted alkyl, cycloalkyl, benzyl, substituted benzyl, or aryl;

4. The compounds according to claim 2, wherein R<sub>1</sub> and R<sub>2</sub> are both H.

5. The compounds according to claim 4, wherein R<sub>5</sub> is independently selected from the group consisting of -H, alkyl, alkenyl, alkynyl, cycloalkyl, heterocycloalkyl, halogenated alkyl, halogenated alkenyl, halogenated alkynyl, halogenated cycloalkyl,  
 25 halogenated heterocycloalkyl, limited substituted alkyl, limited substituted alkenyl, limited substituted alkynyl, substituted cycloalkyl, substituted heterocycloalkyl, -OR<sub>14</sub>, -SR<sub>14</sub>, -F, -Cl, -Br, -I, -NR<sub>14</sub>R<sub>14</sub>, -C(O)R<sub>14</sub>, -C(O)NR<sub>14</sub>R<sub>14</sub>, -CN, -NR<sub>8</sub>C(O)R<sub>11</sub>, -S(O)<sub>2</sub>NR<sub>8</sub>R<sub>8</sub>, -OS(O)<sub>2</sub>R<sub>11</sub>, -S(O)<sub>2</sub>R<sub>14</sub>, -NR<sub>8</sub>S(O)<sub>2</sub>R<sub>8</sub>, -N(H)C(O)N(H)R<sub>8</sub>, and -NO<sub>2</sub>.

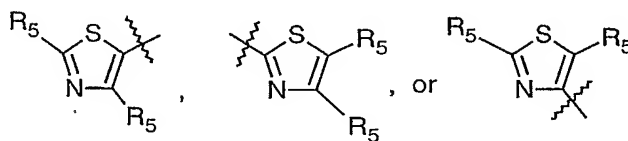
6. The compounds according to claim 5, wherein R<sub>3</sub> is -H, alkyl, cycloalkyl,  
 30 heterocycloalkyl, halogenated alkyl, halogenated cycloalkyl, halogenated heterocycloalkyl, limited substituted alkyl, substituted cycloalkyl, or substituted heterocycloalkyl.

7. The compounds according to claim 4, wherein W is selected from the following:

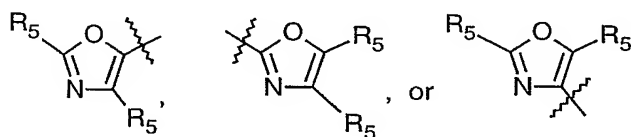


provided that at least one R<sub>5</sub> is not H.

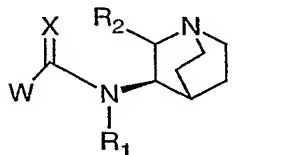
8. The compounds according to claim 4, wherein W is selected from the following:



9. The compounds according to claim 4, wherein W is selected from the following:



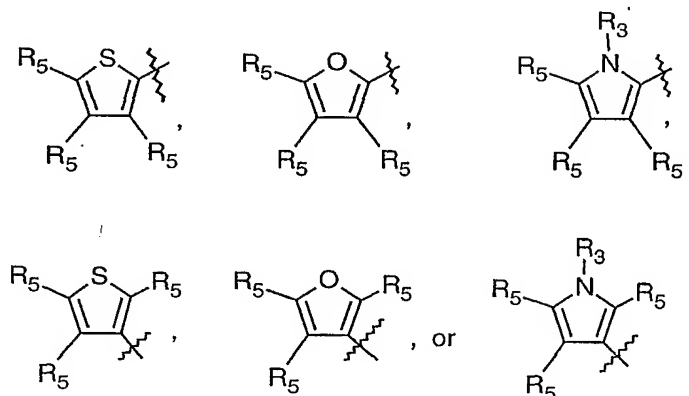
10. The compounds according to claim 1, wherein the compounds are the R stereoisomer at the C3 position of the quinuclidine of Formula I:



or pharmaceutically acceptable salts thereof.

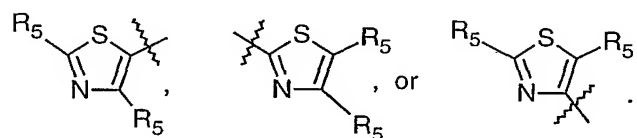
11. The compounds according to claim 10, wherein X is O.
12. The compounds according to claim 11, wherein R<sub>2</sub> is alkyl, halogenated alkyl, substituted alkyl, cycloalkyl, benzyl, substituted benzyl, or aryl.
13. The compounds according to claim 12, wherein R<sub>2</sub> is alkyl, cycloalkyl, or aryl.
14. The compounds according to claim 11, wherein R<sub>1</sub> and R<sub>2</sub> are both H.

15. The compounds according to claim 14, wherein W is selected from the following:

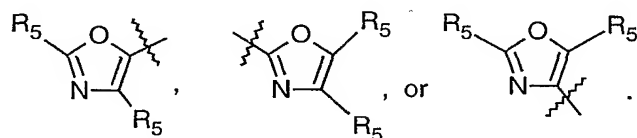


provided that at least one R<sub>5</sub> is not H.

- 5 16. The compounds according to claim 14, wherein W is selected from the following:



17. The compounds according to claim 14, wherein W is selected from the following:



10

18. The compounds according to claim 10, wherein the compounds are selected from the group consisting of:

- 15 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-thiophene-2-carboxamide;  
20 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2,3'-bithiophene-5-carboxamide;

- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-thiophene-2-carboxamide;
- 5 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-phenyl-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-benzyloxyphenyl)-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-benzyloxyphenyl)-thiophene-2-carboxamide;
- 10 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoro-4-benzyloxyphenyl)-thiophene-2-carboxamide;
- 5-(2-aminophenyl)-N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-pyridin-3-yl-thiophene-2-carboxamide;
- 15 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5'-methyl-2,2'-bithiophene-5-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5'-chloro-2,2'-bithiophene-5-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-nitro-thiophene-2-carboxamide;
- 20 5-(aminomethyl)-N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-cyano-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-methoxy-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-2-yl)-thiophene-2-
- 25 carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-[2,2']bithiophenyl-5-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(methylsulfanyl)-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-thiophene-2-carboxamide;
- 30 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-acetyl-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-methyl-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-thiophene-2-carboxamide;

- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylsulfanyl)-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(acetylamino)-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-trifluoromethyl-furan-2-carboxamide;
- 5 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methyl-5-trifluoromethyl-2H-pyrazole-3-yl)-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylthiazol-4-yl)-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(3-chlorophenyl)-vinyl]-thiophene-10 2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenylsulfanyl)-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl-sulfanyl)-thiophene-2-carboxamide;
- 15 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl-sulfanyl)-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chloro-4-fluoro-phenylsulfanyl)-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3-dichlorophenyl-sulfanyl)-20 thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4,5-trichlorophenyl-sulfanyl)-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl-sulfanyl)-thiophene-2-carboxamide;
- 25 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-thiophene-2-carboxamide;
- 30 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoro-4-hydroxyphenyl)-thiophene-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(phenylsulfanyl)-1,3-thiazole-5-carboxamide;

- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[(4-chlorophenyl)-sulfanyl]-1,3-thiazole-5-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-phenoxy-1,3-thiazole-5-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[(4-fluorophenyl)-sulfanyl]-1,3-thiazole-5-carboxamide;
- 5 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(methylsulfanyl)-1,3-thiazole-5-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-chlorophenoxy)-1,3-thiazole-5-carboxamide;
- 10 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-phenyl-1,3-thiazole-5-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2,4-dimethyl-1,3-thiazole-5-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-fluorophenyl)-1,3-thiazole-5-carboxamide;
- 15 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-fluorophenyl)-1,3-thiazole-5-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-fluorophenyl)-1,3-thiazole-5-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-hydroxyphenyl)-1,3-thiazole-5-carboxamide;
- 20 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-methylphenyl)-1,3-thiazole-5-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[4-(benzyloxy)phenyl]-1,3-thiazole-5-carboxamide;
- 25 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-pyridin-3-yl-1,3-thiazole-4-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-2-phenyl-1,3-thiazole-5-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-chlorophenyl)-4-methyl-1,3-thiazole-5-carboxamide;
- 30 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-2-pyridin-2-yl-1,3-thiazole-5-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-2-pyridin-4-yl-1,3-thiazole-5-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(methylamino)-1,3-thiazole-5-carboxamide;

5 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylsulfanyl)-1,3,4-thiadiazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-1,3,4-thiadiazole-2-carboxamide;

10 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-chlorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenoxy)-1,3,4-thiadiazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-1,3,4-thiadiazole-2-carboxamide;

15 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-fluorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-chlorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide;

20 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-fluorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-chlorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1H-1,2,4-triazole-3-carboxamide;

25 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1,3,4-oxadiazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(methylthio)-1,3,4-oxadiazole-2-carboxamide;

30 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1,3-oxazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-1,3-oxazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-1,3-oxazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-1,3-oxazole-2-carboxamide;

5 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-1,3-oxazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1,3-oxazole-2-carboxamide;

10 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1,3-oxazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-1,3-oxazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-1,3-oxazole-2-carboxamide;

15 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-1,3-oxazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-5-phenyl-1,3-thiazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1,3-thiazole-2-carboxamide;

20 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-1,3-thiazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-1,3-thiazole-2-carboxamide;

25 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-1,3-thiazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-1,3-thiazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1,3-thiazole-2-carboxamide;

30 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1,3-thiazole-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-1,3-thiazole-2-carboxamide;

- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-methyl-1,3-thiazole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1,3-thiazole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-1,3-thiazole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1,3-thiazole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-thien-2-yl-1,3-thiazole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(trifluoromethyl)-phenyl]-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(trifluoromethyl)-phenyl]-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,5-difluorophenyl)-furan-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-furan-2-carboxamide;

- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(trifluoromethoxy)phenyl]-furan-2-carboxamide;
- 5 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-chloro-5-(trifluoromethyl)phenyl]-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoro-3-methylphenyl)-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-furan-2-carboxamide;
- 10 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-thien-2-yl-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-thien-3-yl-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-nitro-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4,5-dimethyl-furan-2-carboxamide;
- 15 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloro-2-nitrophenyl)-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methyl-2-nitrophenyl)-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3-difluorophenyl)-furan-2-
- 20 carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenyl)-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-furan-2-carboxamide;
- 25 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxy-phenyl)-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-furan-2-
- 30 carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-furan-2-carboxamide;

- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(trifluoromethoxy)phenyl]-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(trifluoromethoxy)phenyl]-furan-2-carboxamide;
- 5 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-tert-butylphenyl)-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(1-benzothien-2-yl)-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-quinolin-3-yl-furan-2-carboxamide;
- 10 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-ethylphenyl)-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-isopropylphenyl)-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoro-4-methoxyphenyl)-furan-2-carboxamide;
- 15 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(1-benzofuran-2-yl)-furan-2-carboxamide;
- 5-(2-aminophenyl)-N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-furan-2-carboxamide;
- 5-(4-aminophenyl)-N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-furan-2-
- 20 carboxamide;
- 5-(2-amino-4-methylphenyl)-N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylethynyl)-furan-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-furan-2-carboxamide;
- 25 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-1-methyl-5-phenyl-1H-pyrrole-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1,3-oxazole-2-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-3-phenyl-1,2,4-oxadiazole-5-
- carboxamide;
- 30 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-phenyl-1,3-oxazole-5-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-phenyl-1,3-oxazole-4-carboxamide;
- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenylisoxazole-3-carboxamide;
- 5-bromo-N-(2-methyl-1-azabicyclo[2.2.2]oct-3-yl)-thiophene-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-thiophene-2-carbothioamide; and pharmaceutically acceptable salts thereof.

19. The compounds according to claim 10, wherein anyone of or combination of the compounds including:

- 5 N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3'-bithiophene-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-benzyloxyphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-benzyloxyphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoro-4-benzyloxyphenyl)-thiophene-2-carboxamide; 5-(2-aminophenyl)-N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-pyridin-3-yl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5'-methyl-2,2'-bithiophene-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5'-chloro-2,2'-bithiophene-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-nitro-thiophene-2-carboxamide; 5-(aminomethyl)-N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-cyano-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-methoxy-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-[2,2']bithiophenyl-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(methylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-acetyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-methyl-thiophene-2-

carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(acetylamino)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-trifluoromethyl-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methyl-5-trifluoromethyl-2H-pyrazole-3-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylthiazol-4-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(3-chlorophenyl)-vinyl]-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl-sulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl-sulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chloro-4-fluoro-phenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3-dichlorophenyl-sulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4,5-trichlorophenyl-sulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl-sulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoro-4-hydroxyphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(phenylsulfanyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[(4-chlorophenyl)-sulfanyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-phenoxy-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[(4-fluorophenyl)-sulfanyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(methylsulfanyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-chlorophenoxy)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-phenyl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2,4-

dimethyl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-fluorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-fluorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-fluorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-hydroxyphenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-methylphenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[4-(benzyloxy)phenyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-pyridin-3-yl-1,3-thiazole-4-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-2-phenyl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-chlorophenyl)-4-methyl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-2-pyridin-2-yl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-2-pyridin-4-yl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(methylamino)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylsulfanyl)-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-chlorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenoxy)-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-fluorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-chlorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-fluorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-chlorophenyl)-sulfanyl]-1,3,4-thiadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1H-1,2,4-triazole-3-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1,3,4-oxadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(methylthio)-1,3,4-oxadiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-

methoxyphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-5-phenyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-methyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-thien-2-yl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-

bromophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(trifluoromethyl)-phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(trifluoromethyl)-phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,5-difluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(trifluoromethoxy)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-chloro-5-(trifluoromethyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoro-3-methylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-thien-2-yl-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-thien-3-yl-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-nitro-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4,5-dimethyl-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloro-2-nitrophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methyl-2-nitrophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3-difluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-

(2-methylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(trifluoromethoxy)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(trifluoromethoxy)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-tert-butylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(1-benzothien-2-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-quinolin-3-yl-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-ethylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-isopropylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoro-4-methoxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(1-benzofuran-2-yl)-furan-2-carboxamide; 5-(2-aminophenyl)-N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-furan-2-carboxamide; 5-(4-aminophenyl)-N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-furan-2-carboxamide; 5-(2-amino-4-methylphenyl)-N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylethynyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-1-methyl-5-phenyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-3-phenyl-1,2,4-oxadiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-phenyl-1,3-oxazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-phenyl-1,3-oxazole-4-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenylisoxazole-3-carboxamide; 5-bromo-N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-thiophene-2-carbothioamide;

30 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetamidophenyl)-thiophene-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetamidophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetamidophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-difluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-difluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-difluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-carbamoylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-carbamoylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-carbamoylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-sulfamoylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-sulfamoylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-sulfamoylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-

trifluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-difluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-difluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-difluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-carbamoylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-carbamoylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-carbamoylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-sulfamoylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-sulfamoylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-sulfamoylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetamidophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetamidophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetamidophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoroacetamidophenoxy)-

thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-difluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-difluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-difluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-carbamoylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-carbamoylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-carbamoylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-sulfamoylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-sulfamoylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-sulfamoylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-thiophene-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenoxy)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenoxy)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenoxy)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-thiophene-2-carboxamide;  
5 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-thiophene-2-carboxamide;  
10 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethylphenyl)-thiophene-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethylphenyl)-  
thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
trifluoromethylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(2-acetylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
15 yl]-5-(3-acetylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(4-acetylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(2-cyanophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(3-cyanophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(4-cyanophenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
20 yl]-5-(2-morpholin-4-yl-phenyl)-thiophene-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3-morpholin-4-yl-phenyl)-thiophene-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-morpholin-4-yl-phenyl)-thiophene-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenoxy)-  
thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
25 methoxyphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-  
(4-methoxyphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(2-methylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-  
3-yl]-5-(3-methylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenoxy)-thiophene-2-carboxamide; N-[(3R)-  
30 1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethylphenoxy)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethylphenoxy)-thiophene-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethylphenoxy)-  
thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-

acetylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetylphenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-  
5 (3-cyanophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-morpholin-4-yl-phenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-morpholin-4-yl-phenoxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-morpholin-4-yl-phenoxy)-thiophene-2-  
10 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenylsulfanyl)-thiophene-2-carboxamide; N-  
15 [(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethylphenylsulfanyl)-thiophene-2-  
20 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetylphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetylphenylsulfanyl)-  
25 thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-morpholin-4-yl-phenylsulfanyl)-thiophene-2-  
30 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-morpholin-4-yl-phenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-morpholin-4-yl-phenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-yl)-thiophene-2-carboxamide; N-[(3R)-1-

azabicyclo[2.2.2]oct-3-yl]-5-(2-methylpyridin-4-yl)-thiophene-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methylpyridin-3-yl)-thiophene-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylpyridin-2-yl)-  
thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
5 methoxypyridin-4-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(6-methoxypyridin-3-yl)-thiophene-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxypyridin-2-yl)-thiophene-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-yl)-thiophene-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-yl)-thiophene-2-  
10 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-yl)-  
thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-  
3-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-  
chloropyridin-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-  
5-(4-chloropyridin-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-  
15 3-yl]-5-(pyridin-4-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-  
3-yl]-5-(pyridin-3-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-  
3-yl]-5-(pyridin-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-  
3-yl]-5-(2-methylpyridin-4-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(6-methylpyridin-3-yloxy)-thiophene-2-carboxamide; N-  
20 [(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylpyridin-2-yloxy)-thiophene-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxypyridin-4-yloxy)-  
thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-  
methoxypyridin-3-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-  
3-yl]-5-(5-methoxypyridin-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-  
25 azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-yloxy)-thiophene-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-yloxy)-thiophene-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-yloxy)-  
thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-  
4-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-  
30 chloropyridin-3-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(5-chloropyridin-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-3-yloxy)-thiophene-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-2-yloxy)-thiophene-2-

carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylpyridin-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methylpyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylpyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxypyridin-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methoxypyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxypyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiophen-4-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiophen-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiophen-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiophen-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(furan-4-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylfuran-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyfuran-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorofuran-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-2-yl)-thiophene-2-

carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyloxazol-2-yl)-  
thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyoxazol-  
2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-  
chlorooxazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-  
5-(oxazol-5-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
methyloxazol-5-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-  
5-(2-methoxyoxazol-5-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-  
3-yl]-5-(2-chlorooxazol-5-yl)-thiophene-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiazol-2-yl)-thiophene-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiazol-2-yl)-thiophene-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiazol-2-yl)-  
thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-5-yl)-  
thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylthiazol-  
5-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
methoxythiazol-5-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(2-chlorothiazol-5-yl)-thiophene-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(4-methylthiazol-2-yl)-thiophene-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxythiazol-2-yl)-thiophene-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorothiazol-2-yl)-  
thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methyloxazol-  
2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
methoxyoxazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(4-chlorooxazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]thiadiazol-2-yl)-thiophene-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]thiadiazol-2-yl)-thiophene-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]thiadiazol-2-  
yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-  
chloro[1,3,4]thiadiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]oxadiazol-2-yl)-thiophene-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]oxadiazol-2-yl)-thiophene-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]oxadiazol-2-  
yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-

chloro[1,3,4]oxadiazol-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiophen-4-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiophen-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiophen-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiophen-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(furan-4-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylfuran-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyfuran-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorofuran-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyloxazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyoxazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorooxazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-5-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methyloxazol-5-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyoxazol-5-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorooxazol-5-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-5-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylthiazol-5-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxythiazol-5-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorothiazol-5-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylthiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxythiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorothiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methyloxazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-

methoxyoxazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorooxazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]thiadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]thiadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]thiadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]thiadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]oxadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]oxadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]oxadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]oxadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiophen-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiophen-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiophen-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiophen-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(furan-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylfuran-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyfuran-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorofuran-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyloxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyoxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorooxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methyloxazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyoxazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorooxazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-

[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylthiazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxythiazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorothiazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylthiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxythiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorothiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methyloxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyoxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorooxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]thiadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]thiadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]thiadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]thiadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]oxadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]oxadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]oxadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]oxadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrrole-2-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(isothiazol-3-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(isoxazol-3-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3H-imidazol-4-yl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-(4-hydroxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-(3-acetamidophenylsulfanyl)-thiophene-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-(2-methanesulfonylamino-phenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-(pyridin-2-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-(2-methylpyridin-4-yloxy)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-(4-trifluoromethylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-(2-acetylphenyl)-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-phenyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-5-phenyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-cyano-5-phenyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-methoxy-5-phenyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-phenyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-methyl-4-phenyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-cyano-4-phenyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-methoxy-4-phenyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-bromo-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-methylsulfanyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-methyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-cyano-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-bromo-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-methylsulfanyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-methyl-thiophene-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-cyano-thiophene-2-carboxamide;

N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methanesulfonylamino-phenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-

azabicyclo[2.2.2]oct-3-yl]-5-(3-methanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-difluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-difluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-difluoroacetamidophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-carbamoylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-carbamoylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-carbamoylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-sulfamoylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-sulfamoylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-sulfamoylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-

- trifluoromethanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
trifluoromethanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
5 trifluoromethanesulfonylaminophenylsulfanyl)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
difluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3-difluoroacetamidophenylsulfanyl)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
10 difluoroacetamidophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-carbamoylphenylsulfanyl)-thiophene-2-carboxamide;  
N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-carbamoylphenylsulfanyl)-  
thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
carbamoylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
15 azabicyclo[2.2.2]oct-3-yl]-5-(2-sulfamoylphenylsulfanyl)-thiophene-2-carboxamide;  
N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-sulfamoylphenylsulfanyl)-  
thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
sulfamoylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenylsulfanyl)-thiophene-2-carboxamide; N-  
20 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenylsulfanyl)-  
thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
hydroxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenylsulfanyl)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenylsulfanyl)-thiophene-  
25 2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
aminophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-acetamidophenoxy)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetamidophenoxy)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
30 acetamidophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoroacetamidophenoxy)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
trifluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-

azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethanesulfonylaminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-difluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-difluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-difluoroacetamidophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-carbamoylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-carbamoylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-carbamoylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-sulfamoylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-sulfamoylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-sulfamoylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenoxy)-thiophene-2-

carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-morpholin-4-yl-phenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-morpholin-4-yl-phenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-morpholin-4-yl-phenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-

azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenoxy)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenoxy)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
methylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
5 azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethylphenoxy)-thiophene-2-carboxamide;  
N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-trifluoromethylphenoxy)-  
thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
trifluoromethylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-acetylphenoxy)-thiophene-2-carboxamide; N-  
10 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetylphenoxy)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
acetylphenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenoxy)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenoxy)-thiophene-2-  
15 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
cyanophenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-morpholin-4-yl-phenoxy)-thiophene-2-carboxamide;  
N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-morpholin-4-yl-phenoxy)-  
thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
20 morpholin-4-yl-phenoxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenylsulfanyl)-thiophene-2-carboxamide;  
N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenylsulfanyl)-  
thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
methoxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
25 azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenylsulfanyl)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenylsulfanyl)-  
thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
methylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-trifluoromethylphenylsulfanyl)-thiophene-2-  
30 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
trifluoromethylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(4-trifluoromethylphenylsulfanyl)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-

acetylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetylphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-morpholin-4-yl-phenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-morpholin-4-yl-phenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-morpholin-4-yl-phenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylpyridin-4-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methylpyridin-3-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylpyridin-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxypyridin-4-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methoxypyridin-3-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxypyridin-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-3-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-3-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylpyridin-4-yloxy)-

thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methylpyridin-3-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylpyridin-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxypyridin-4-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methoxypyridin-3-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxypyridin-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-3-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylpyridin-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methylpyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylpyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxypyridin-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-methoxypyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxypyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-

3-yl]-5-(6-chloropyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-3-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiophen-4-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiophen-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiophen-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiophen-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(furan-4-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylfuran-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyfuran-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorofuran-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyloxazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyoxazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorooxazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-5-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methyloxazol-5-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyoxazol-5-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorooxazol-5-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-

azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiazol-2-yl)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiazol-2-yl)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-5-yl)-  
thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
5 methylthiazol-5-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxythiazol-5-yl)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorothiazol-5-yl)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylthiazol-  
2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-  
10 (4-methoxythiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorothiazol-2-yl)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methyloxazol-2-yl)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
methoxyoxazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
15 azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorooxazol-2-yl)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]thiadiazol-2-yl)-thiophene-  
2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-  
methyl[1,3,4]thiadiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]thiadiazol-2-yl)-thiophene-2-  
20 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-  
chloro[1,3,4]thiadiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]oxadiazol-2-yl)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]oxadiazol-2-yl)-  
thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-  
25 methoxy[1,3,4]oxadiazol-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]oxadiazol-2-yl)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiophen-4-  
yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-  
5-(5-methylthiophen-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
30 azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiophen-2-yloxy)-thiophene-2-carboxamide;  
N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiophen-2-yloxy)-  
thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-  
(furan-4-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-

azabicyclo[2.2.2]oct-3-yl]-5-(5-methylfuran-2-yloxy)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyfuran-2-yloxy)-  
thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-  
chlorofuran-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
5 azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-  
2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyloxazol-2-yloxy)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-  
methoxyoxazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorooxazol-2-yloxy)-thiophene-2-carboxamide; N-  
10 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-5-yloxy)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methyloxazol-  
5-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(2-methoxyoxazol-5-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorooxazol-5-yloxy)-thiophene-2-carboxamide; N-  
15 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-2-yloxy)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiazol-  
2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(5-methoxythiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiazol-2-yloxy)-thiophene-2-carboxamide; N-  
20 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-5-yloxy)-thiophene-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylthiazol-  
5-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(2-methoxythiazol-5-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorothiazol-5-yloxy)-thiophene-2-carboxamide; N-  
25 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylthiazol-2-yloxy)-  
thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
methoxythiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorothiazol-2-yloxy)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methyloxazol-2-yloxy)-  
30 thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
methoxyoxazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorooxazol-2-yloxy)-thiophene-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]thiadiazol-2-yloxy)-

thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]thiadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]thiadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]thiadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]oxadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]oxadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]oxadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]oxadiazol-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiophen-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiophen-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiophen-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiophen-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(furan-4-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylfuran-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyfuran-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorofuran-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyloxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxyoxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorooxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(oxazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methyloxazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyoxazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorooxazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-

[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methylthiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxythiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chlorothiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(thiazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylthiazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxythiazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorothiazol-5-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylthiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxythiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorothiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methyloxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyoxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorooxazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]thiadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]thiadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]thiadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]thiadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-([1,3,4]oxadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methyl[1,3,4]oxadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-methoxy[1,3,4]oxadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloro[1,3,4]oxadiazol-2-ylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrrole-2-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(isothiazol-3-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(isoxazol-3-yl)-thiophene-2-carboxamide;

N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3H-imidazol-4-yl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-(4-hydroxyphenylsulfanyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-(3-acetamidophenylsulfanyl)-thiophene-2-carboxamide;

5 N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-(2-methanesulfonylaminophenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-(pyridin-2-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-(2-methylpyridin-4-yloxy)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-(4-trifluoromethylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-(2-acetylphenyl)-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-5-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-cyano-5-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-methoxy-5-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-methyl-4-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-cyano-4-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-methoxy-4-phenyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-bromo-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-methylsulfanyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-methyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-chloro-5-cyano-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-bromo-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-methylsulfanyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-methyl-thiophene-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-4-cyano-thiophene-2-carboxamide;

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N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3-dichlorophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,5-dichlorophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,6-dichlorophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-amino-2-fluorophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-amino-2-fluorophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-amino-2-chlorophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-amino-2-chlorophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoro-4-methylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoro-3-methoxyphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoro-4-methoxyphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-[(methylsulfonyl)amino]phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-[(methylsulfonyl)amino]phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-[(methylsulfonyl)amino]phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-[(trifluoromethyl)sulfonyl]amino)phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-[(trifluoromethyl)sulfonyl]amino)phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-[(trifluoromethyl)sulfonyl]amino)phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-[(phenylsulfonyl)amino]phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-[(phenylsulfonyl)amino]phenyl)-furan-2-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-[(phenylsulfonyl)amino]phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-[(methylamino)carbonyl]phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-[(methylamino)carbonyl]phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-[(methylamino)carbonyl]phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-[(methylamino)sulfonyl]phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-[(methylamino)sulfonyl]phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-[(methylamino)sulfonyl]phenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(methylamino)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(methylamino)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(methylamino)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(ethylamino)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(ethylamino)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(ethylamino)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-acetylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(methylthio)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(methylthio)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(methylthio)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(phenylthio)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(phenylthio)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(phenylthio)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-phenoxyphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-phenoxyphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-phenoxyphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-anilinophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-anilinophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-anilinophenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylthio)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-

fluorophenyl)thio]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-fluorophenyl)thio]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-fluorophenyl)thio]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-chlorophenyl)thio]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-chlorophenyl)thio]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-chlorophenyl)thio]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenoxy)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenoxy)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenoxy)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenoxy)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenoxy)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-pyridin-2-yl-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-pyridin-3-yl-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-pyridin-4-yl-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyridin-2-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-2-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-2-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-2-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-3-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-3-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-3-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-3-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-4-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyridin-4-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chloropyridin-2-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-2-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-2-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-3-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-3-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-3-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-yl)-furan-2-

carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chloropyridin-4-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-piperidin-1-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-piperidin-1-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-piperidin-1-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-morpholin-4-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-morpholin-4-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-morpholin-4-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-pyrrolidin-1-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-pyrrolidin-1-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-pyrrolidin-1-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(1H-pyrrol-2-yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(1H-pyrrol-2-yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1H-pyrrol-2-yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(2-furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(3-furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(5-methyl-2-furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(2-furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(3-furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(5-methyl-2-furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(2-furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(3-furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(5-methyl-2-furyl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(1,3-thiazol-2-yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(1,3-thiazol-2-yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1,3-thiazol-2-yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(1,3-oxazol-2-yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(1,3-oxazol-2-yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1,3-oxazol-2-yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-isothiazol-5-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-isothiazol-5-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-

5-(4-isothiazol-5-ylphenyl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1H-indol-2-yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1H-indol-3-yl)phenyl]-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(1H-indol-5-yl)-furan-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(1H-indol-6-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3-dichlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,5-dichlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,6-dichlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-amino-2-fluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-amino-2-fluorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-amino-2-chlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-amino-2-chlorophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoro-4-methylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoro-3-methoxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoro-4-methoxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-[(methylsulfonyl)amino]phenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-[(methylsulfonyl)amino]phenyl)-furan-2-

carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
[(methylsulfonyl)amino]phenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-[2-([(trifluoromethyl)sulfonyl]amino)phenyl]-furan-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-  
5 [(trifluoromethyl)sulfonyl]amino)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-  
methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-([(trifluoromethyl)sulfonyl]amino)phenyl)-  
furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
[(phenylsulfonyl)amino]phenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3-[(phenylsulfonyl)amino]phenyl)-furan-2-carboxamide;  
10 N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
[(phenylsulfonyl)amino]phenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-[(methylamino)carbonyl]phenyl)-furan-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
[(methylamino)carbonyl]phenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
15 azabicyclo[2.2.2]oct-3-yl]-5-(4-[(methylamino)carbonyl]phenyl)-furan-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
[(methylamino)sulfonyl]phenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3-[(methylamino)sulfonyl]phenyl)-furan-2-carboxamide;  
N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
20 [(methylamino)sulfonyl]phenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-[2-(methylamino)phenyl]-furan-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(methylamino)phenyl]-furan-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-  
(methylamino)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
25 azabicyclo[2.2.2]oct-3-yl]-5-[2-(ethylamino)phenyl]-furan-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(ethylamino)phenyl]-furan-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-  
(ethylamino)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-acetylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-  
30 methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-acetylphenyl)-furan-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-acetylphenyl)-furan-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-  
(methylthio)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-

azabicyclo[2.2.2]oct-3-yl]-5-[3-(methylthio)phenyl]-furan-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(methylthio)phenyl]-furan-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-  
(phenylthio)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
5 azabicyclo[2.2.2]oct-3-yl]-5-[3-(phenylthio)phenyl]-furan-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(phenylthio)phenyl]-furan-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
phenoxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-  
3-yl]-5-(3-phenoxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
10 azabicyclo[2.2.2]oct-3-yl]-5-(4-phenoxyphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-  
methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-anilinophenyl)-furan-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-anilinophenyl)-furan-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-anilinophenyl)-  
furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-  
15 (phenylthio)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-  
yl]-5-[(2-fluorophenyl)thio]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-[(3-fluorophenyl)thio]-furan-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-fluorophenyl)thio]-furan-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-  
20 chlorophenyl)thio]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-[(3-chlorophenyl)thio]-furan-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-chlorophenyl)thio]-furan-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
fluorophenoxy)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-  
25 yl]-5-(3-fluorophenoxy)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenoxy)-furan-2-carboxamide; N-[(2S,3R)-2-  
methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenoxy)-furan-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenoxy)-furan-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
30 chlorophenoxy)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-  
3-yl]-5-pyridin-2-yl-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-pyridin-3-yl-furan-2-carboxamide; N-[(2S,3R)-2-methyl-  
1-azabicyclo[2.2.2]oct-3-yl]-5-pyridin-4-yl-furan-2-carboxamide; N-[(2S,3R)-2-

methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyridin-2-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-2-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-2-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-2-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-3-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-3-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-3-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-3-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-4-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyridin-4-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chloropyridin-2-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-2-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-2-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-2-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-3-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chloropyridin-3-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-chloropyridin-3-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-chloropyridin-3-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chloropyridin-4-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chloropyridin-4-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-piperidin-1-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-piperidin-1-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-piperidin-1-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-morpholin-4-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-morpholin-4-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-morpholin-4-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-pyrrolidin-1-

ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-pyrrolidin-1-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-pyrrolidin-1-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(1H-pyrrol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(1H-pyrrol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1H-pyrrol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(2-furyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(3-furyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(5-methyl-2-furyl)phenyl]-furan-2-carboxamid; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(2-furyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(3-furyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(5-methyl-2-furyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(2-furyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(3-furyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(5-methyl-2-furyl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(1,3-thiazol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(1,3-thiazol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1,3-thiazol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(1,3-oxazol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(1,3-oxazol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1,3-oxazol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-isothiazol-5-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-isothiazol-5-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-isothiazol-5-ylphenyl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1H-indol-2-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(1H-indol-3-yl)phenyl]-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-

5-(1H-indol-5-yl)-furan-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(1H-indol-6-yl)-furan-2-carboxamide;

- N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2,3-difluorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2,4-difluorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2,5-difluorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2,6-difluorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3,4-difluorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3,5-difluorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-chlorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-chlorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-chlorophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-bromophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-bromophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-bromophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-cyanophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-cyanophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-cyanophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-nitrophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-nitrophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-nitrophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-methylphenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-methylphenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-aminophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-aminophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-aminophenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[2-(methylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[3-(methylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[4-(methylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[2-(acetylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[3-(acetylamino)phenyl]-1,3-thiazole-5-carboxamide; N-

[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[4-(acetylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-[(trifluoroacetyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-[(trifluoroacetyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-[(trifluoroacetyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-[(methylsulfonyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-[(methylsulfonyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-[(methylsulfonyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-hydroxyphenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-hydroxyphenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-methoxyphenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-methoxyphenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-methoxyphenyl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[2-(trifluoromethoxy)phenyl]-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[3-(trifluoromethoxy)phenyl]-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-[4-(trifluoromethoxy)phenyl]-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-pyridin-2-yl-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-pyridin-3-yl-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-pyridin-4-yl-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(6-fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(5-fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(6-fluoropyridin-3-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(5-fluoropyridin-3-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-fluoropyridin-3-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-fluoropyridin-3-yl)-1,3-

- thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-fluoropyridin-4-yl)-1,3-thiazole-5-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-fluoropyridin-4-yl)-1,3-thiazole-5-carboxamide;
- 5 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(methylthio)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylthio)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-fluorophenyl)thio]-1,3-thiazole-2-carboxamide; N-
- 10 [(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-fluorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-fluorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-chlorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-chlorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(3R)-1-
- 15 azabicyclo[2.2.2]oct-3-yl]-5-[(4-chlorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenoxy)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenoxy)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenoxy)-1,3-thiazole-2-
- 20 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenoxy)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenoxy)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-thien-3-yl-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-
- 25 (2,4-difluorophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-1,3-thiazole-2-
- 30 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-

(4-cyanophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-(acetylamino)phenyl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(acetylamino)phenyl]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetylamino)phenyl]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-4-yl)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-methoxyphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-methoxyphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-methoxyphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(3R)-1-

azabicyclo[2.2.2]oct-3-yl]-5-[(2-methylphenyl)thio]-1,3-thiazole-2-carboxamide; N-  
 [(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-methylphenyl)thio]-1,3-thiazole-2-  
 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-methylphenyl)thio]-1,3-  
 thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-  
 5 (acetylamino)phenyl)thio]-1,3-thiazole-2-carboxamide; N-[(3R)-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-[(4-aminophenyl)thio]-1,3-thiazole-2-carboxamide; N-  
 [(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-hydroxyphenyl)thio]-1,3-thiazole-2-  
 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenoxy)-1,3-  
 thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenoxy)-  
 10 1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
 methylphenoxy)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-  
 (2-methoxyphenoxy)-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
 yl]-5-(3-methoxyphenoxy)-1,3-thiazole-2-carboxamide; N-[(3R)-1-  
 azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenoxy)-1,3-thiazole-2-carboxamide; N-  
 15 [(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenoxy)-1,3-thiazole-2-  
 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenoxy)-1,3-  
 thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-  
 (acetylamino)phenoxy]-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-  
 3-yl]-5-(2-fluorophenyl)-4-methyl-1,3-thiazole-2-carboxamide; N-[(3R)-1-  
 20 azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-4-methyl-1,3-thiazole-2-carboxamide;  
 N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-4-methyl-1,3-thiazole-2-  
 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-4-methyl-1,3-  
 thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-  
 4-methyl-1,3-thiazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
 25 methoxyphenyl)-4-methyl-1,3-thiazole-2-carboxamide; N-[(3R)-1-  
 azabicyclo[2.2.2]oct-3-yl]-4-methyl-5-thien-2-yl-1,3-thiazole-2-carboxamide,  
  
 N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2,3-difluorophenyl)-1,3-  
 thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2,4-  
 30 difluorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-  
 azabicyclo[2.2.2]oct-3-yl]-2-(2,5-difluorophenyl)-1,3-thiazole-5-carboxamide; N-  
 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2,6-difluorophenyl)-1,3-thiazole-5-  
 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3,4-

5 difluorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3,5-difluorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-chlorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-chlorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-chlorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-bromophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-bromophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-bromophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-cyanophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-cyanophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-cyanophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-nitrophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-nitrophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-nitrophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-methylphenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-methylphenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-aminophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-aminophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-aminophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[2-(methylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[3-(methylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[4-(methylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[2-(acetylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[3-(acetylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[4-(acetylamino)phenyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-[(trifluoroacetyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-

azabicyclo[2.2.2]oct-3-yl]-2-(3-[(trifluoroacetyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-[(trifluoroacetyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-[(methylsulfonyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-[(methylsulfonyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-[(methylsulfonyl)amino]phenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-hydroxyphenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-hydroxyphenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-methoxyphenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-methoxyphenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-methoxyphenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[2-(trifluoromethoxy)phenyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[3-(trifluoromethoxy)phenyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-[4-(trifluoromethoxy)phenyl]-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-pyridin-2-yl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-pyridin-3-yl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-pyridin-4-yl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(6-fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(5-fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-fluoropyridin-2-yl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(6-fluoropyridin-3-yl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(5-fluoropyridin-3-yl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-fluoropyridin-3-yl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-

fluoropyridin-3-yl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-fluoropyridin-4-yl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-fluoropyridin-4-yl)-1,3-thiazole-5-carboxamide;

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N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(methylthio)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-chloro-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-1,3-thiazole-2-carboxamide, N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(phenylthio)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-fluorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-fluorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-fluorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-chlorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-chlorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-chlorophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-thien-3-yl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-1,3-thiazole-2-

carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-1,3-thiazole-2-carboxamide; N-  
5 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-1,3-thiazole-2-carboxamide; N-  
10 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-1,3-thiazole-2-carboxamide; N-  
15 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenyl)-1,3-thiazole-2-carboxamide; N-  
20 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(acetylamino)phenyl]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(acetylamino)phenyl]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetylamino)phenyl]-1,3-  
25 thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyridin-2-yl)-1,3-  
30 thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-2-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-2-yl)-1,3-thiazole-

2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-3-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-4-yl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-methoxyphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-methoxyphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-methoxyphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(2-methylphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(3-methylphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-methylphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-(acetylamino)phenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-aminophenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[(4-hydroxyphenyl)thio]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenoxy)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetylamino)phenoxy]-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-4-methyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-

fluorophenyl)-4-methyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-4-methyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-4-methyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-4-methyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-4-methyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-4-methyl-5-thien-2-yl-1,3-thiazole-2-carboxamide, N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-4-methyl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-thien-2-yl-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-1,3-thiazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-hydroxyphenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-methylphenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-phenyl-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(2-fluorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(3-fluorophenyl)-1,3-thiazole-5-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-2-(4-fluorophenyl)-1,3-thiazole-5-carboxamide;

N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenylsulfanyl-1,3-oxazole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenylsulfanyl)-1,3-oxazole-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenylsulfanyl)-1,3-  
5 oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
fluorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(3,4-difluorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenylsulfanyl)-1,3-oxazole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenylsulfanyl)-1,3-oxazole-2-  
10 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenylsulfanyl)-1,3-  
oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
chlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(4-chlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenylsulfanyl)-1,3-oxazole-2-  
15 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenylsulfanyl)-  
1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-  
dichlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-  
3-yl]-5-(2-fluorophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-  
20 1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenyl)-1,3-oxazole-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-1,3-oxazole-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-1,3-  
oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-  
1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-  
25 dichlorophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-  
(2,4-dichlorophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(3,5-dichlorophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-  
1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-1,3-oxazole-2-carboxamide; N-  
30 [(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-1,3-oxazole-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-1,3-oxazole-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-1,3-oxazole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-1,3-oxazole-2-

carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylaminophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylaminophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylaminophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenyl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(acetylamino)phenyl]-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(acetylamino)phenyl]-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetylamino)phenyl]-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-2-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-3-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyridin-2-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-2-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-2-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-2-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-3-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-3-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-3-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-3-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-4-yl)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-

5-(3-fluorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenoxy)-1,3-oxazole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenoxy)-1,3-oxazole-2-carboxamide;

N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenylsulfanyl-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenylsulfanyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1,3-oxazole-2-

carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylaminophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylaminophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylaminophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenyl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(acetylamino)phenyl]-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-

azabicyclo[2.2.2]oct-3-yl]-5-[3-(acetylamino)phenyl]-1,3-oxazole-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetylamino)phenyl]-1,3-  
oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-  
(pyridin-2-yl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
5 azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-3-yl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-  
2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyridin-4-yl)-1,3-oxazole-2-carboxamide;  
N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyridin-2-yl)-1,3-  
oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
fluoropyridin-2-yl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
10 azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-2-yl)-1,3-oxazole-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyridin-2-yl)-1,3-oxazole-  
2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
fluoropyridin-3-yl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyridin-3-yl)-1,3-oxazole-2-carboxamide; N-  
15 [(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyridin-3-yl)-1,3-oxazole-  
2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-  
fluoropyridin-3-yl)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-4-yl)-1,3-oxazole-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyridin-3-yl)-1,3-oxazole-  
20 2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenoxy-1,3-  
oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
fluorophenoxy)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenoxy)-1,3-oxazole-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenoxy)-1,3-oxazole-2-  
25 carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-  
difluorophenoxy)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenoxy)-1,3-oxazole-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenoxy)-1,3-oxazole-  
2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
30 chlorophenoxy)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenoxy)-1,3-oxazole-2-carboxamide; N-  
[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenoxy)-1,3-oxazole-2-  
carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-

dichlorophenoxy)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenoxy)-1,3-oxazole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenoxy)-1,3-oxazole-2-carboxamide;

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N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(methylthio)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(methylthio)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,6-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4,6-trifluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-

azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-1-methyl-1H-pyrrole-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1-methyl-  
1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-  
5 methoxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-1-methyl-1H-pyrrole-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-1-methyl-1H-  
pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-1-  
10 methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
methylaminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3-methylaminophenyl)-1-methyl-1H-pyrrole-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylaminophenyl)-1-  
methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-  
15 nitrophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(3-nitrophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenyl)-1-methyl-1H-pyrrole-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenyl)-1-methyl-1H-  
20 pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenyl)-1-  
methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-  
(acetyl amino)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-[3-(acetyl amino)phenyl]-1-methyl-1H-pyrrole-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetyl amino)phenyl]-1-  
25 methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-2-  
yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-  
(pyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(pyridin-4-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyrid-2-yl)-1-methyl-1H-pyrrole-2-  
30 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyrid-2-yl)-1-methyl-  
1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyrid-2-  
yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-  
fluoropyrid-2-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-

azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyrid-4-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyrid-4-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-1H-pyrrole-2-carboxamide; N-

[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-1H-pyrrole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-1H-pyrrole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-1H-pyrrole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylaminophenyl)-1H-pyrrole-2-  
5 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylaminophenyl)-1H-  
pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-  
methylaminophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(2-nitrophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(3-nitrophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
10 yl]-5-(4-nitrophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(2-aminophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(3-aminophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(4-aminophenyl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-[2-(acetylamino)phenyl]-1H-pyrrole-2-carboxamide; N-[(3R)-1-  
15 azabicyclo[2.2.2]oct-3-yl]-5-[3-(acetylamino)phenyl]-1H-pyrrole-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetylamino)phenyl]-1H-pyrrole-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-2-yl)-1H-pyrrole-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-3-yl)-1H-pyrrole-2-  
carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-4-yl)-1H-pyrrole-2-  
20 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyrid-2-yl)-1H-  
pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyrid-2-yl)-  
1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyrid-2-  
yl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-  
fluoropyrid-2-yl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-  
25 (2-fluoropyrid-3-yl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-  
yl]-5-(4-fluoropyrid-3-yl)-1H-pyrrole-2-carboxamide; N-[(3R)-1-  
azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyrid-3-yl)-1H-pyrrole-2-carboxamide; N-  
[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyrid-3-yl)-1H-pyrrole-2-carboxamide;  
N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyrid-4-yl)-1H-pyrrole-2-  
30 carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyrid-4-yl)-1H-  
pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-1H-  
pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-1H-

pyrrole-2-carboxamide; N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-1H-pyrrole-2-carboxamide;

- 5 N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-bromo-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(methylthio)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(methylthio)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-phenyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,3-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,6-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4,6-trifluorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-1-methyl-1H-pyrrole-2-carboxamide;

N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylaminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylaminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylaminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(acetyl amino)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(acetyl amino)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetyl amino)phenyl]-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-2-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-pyridin-4-yl-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyrid-2-yl)-1-methyl-1H-

pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyrid-2-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyrid-2-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyrid-2-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(5-fluoropyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyrid-3-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyrid-4-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyrid-4-yl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-cyanophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-1-methyl-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-difluorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-difluorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-difluorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-chlorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-chlorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-chlorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,5-dichlorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3,4-dichlorophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2,4-dichlorophenyl)-1H-pyrrole-2-carboxamide; N-

[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-bromophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-bromophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-bromophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-hydroxyphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-hydroxyphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-hydroxyphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methoxyphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methoxyphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methoxyphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methylphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylphenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-methyaminophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-methylaminophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-methylaminophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-nitrophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-nitrophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-nitrophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-aminophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-aminophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-aminophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[2-(acetylamino)phenyl]-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[3-(acetylamino)phenyl]-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-[4-(acetylamino)phenyl]-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-2-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-3-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-

azabicyclo[2.2.2]oct-3-yl]-5-(pyrid-4-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyrid-2-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyrid-2-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-  
5 (5-fluoropyrid-2-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyrid-2-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyrid-3-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-fluoropyrid-3-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-  
10 (5-fluoropyrid-3-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(6-fluoropyrid-3-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(2-fluoropyrid-4-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-fluoropyrid-4-yl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-  
15 (2-cyanophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(3-cyanophenyl)-1H-pyrrole-2-carboxamide; N-[(2S,3R)-2-methyl-1-azabicyclo[2.2.2]oct-3-yl]-5-(4-cyanophenyl)-1H-pyrrole-2-carboxamide; or pharmaceutically acceptable salts thereof.

- 20 20. A method for treating a disease or condition in a mammal, wherein the  $\alpha 7$  nicotinic acetylcholine receptor is implicated comprising administering to a mammal a therapeutically effective amount of compound(s) according to claim 1.
21. The method according to claim 20, wherein the disease or condition is schizophrenia, or psychosis.
- 25 22. The method according to claim 21, wherein said compound(s) is(are) administered rectally, topically, orally, sublingually, or parenterally.
23. The method according to claim 21, wherein said compound(s) is(are) administered from about 0.001 to about 100 mg/kg of body weight of said mammal per day.
- 30 24. The method according to claim 21, wherein said compound(s) is(are) administered from about 0.1 to about 50 mg/kg of body weight of said mammal per day.

25. A method for treating a disease or condition in a mammal, wherein the  $\alpha 7$  nicotinic acetylcholine receptor is implicated comprising administering to a mammal a therapeutically effective amount of compound(s) according to claim 5.

26. The method according to claim 25, wherein the disease or condition is  
5 schizophrenia, or psychosis.

27. The method according to claim 26, wherein said compound(s) is(are) administered rectally, topically, orally, sublingually, or parenterally.

28. The method according to claim 26, wherein said compound(s) is(are) administered from about 0.001 to about 100 mg/kg of body weight of said mammal  
10 per day.

29. The method according to claim 26, wherein said compound(s) is(are) administered from about 0.1 to about 50 mg/kg of body weight of said mammal per day.

30. A method for treating a disease or condition in a mammal, wherein the  $\alpha 7$  nicotinic acetylcholine receptor is implicated comprising administering to a mammal a  
15 therapeutically effective amount of compound(s) according to claim 6.

31. The method according to claim 30, wherein the disease or condition is schizophrenia, or psychosis.

32. The method according to claim 31, wherein said compound(s) is(are)  
20 administered rectally, topically, orally, sublingually, or parenterally.

33. The method according to claim 31, wherein said compound(s) is(are) administered from about 0.001 to about 100 mg/kg of body weight of said mammal per day.

34. The method according to claim 31, wherein said compound(s) is(are)  
25 administered from about 0.1 to about 50 mg/kg of body weight of said mammal per day.

35. A method for treating a disease or condition in a mammal, wherein the  $\alpha 7$  nicotinic acetylcholine receptor is implicated comprising administering to a mammal a therapeutically effective amount of compound(s) according to claim 10.

30 36. The method according to claim 35, wherein the disease or condition is schizophrenia, or psychosis.

37. The method according to claim 36, wherein said compound(s) is(are) administered rectally, topically, orally, sublingually, or parenterally.

38. The method according to claim 36, wherein said compound(s) is(are) administered from about 0.001 to about 100 mg/kg of body weight of said mammal per day.
39. The method according to claim 36, wherein said compound(s) is(are) administered from about 0.1 to about 50 mg/kg of body weight of said mammal per day.
40. A method for treating a disease or condition in a mammal, wherein the  $\alpha 7$  nicotinic acetylcholine receptor is implicated comprising administering to a mammal a therapeutically effective amount of compound(s) according to claim 11.
41. The method according to claim 40, wherein the disease or condition is schizophrenia, or psychosis.
42. The method according to claim 41, wherein said compound(s) is(are) administered rectally, topically, orally, sublingually, or parenterally.
43. The method according to claim 41, wherein said compound(s) is(are) administered from about 0.001 to about 100 mg/kg of body weight of said mammal per day.
44. The method according to claim 41, wherein said compound(s) is(are) administered from about 0.1 to about 50 mg/kg of body weight of said mammal per day.
45. A method for treating a disease or condition in a mammal, wherein the  $\alpha 7$  nicotinic acetylcholine receptor is implicated comprising administering to a mammal a therapeutically effective amount of compound(s) according to claim 12.
46. The method according to claim 45, wherein the disease or condition is schizophrenia, or psychosis.
47. The method according to claim 46, wherein said compound(s) is(are) administered rectally, topically, orally, sublingually, or parenterally.
48. The method according to claim 46, wherein said compound(s) is(are) administered from about 0.001 to about 100 mg/kg of body weight of said mammal per day.
49. The method according to claim 46, wherein said compound(s) is(are) administered from about 0.1 to about 50 mg/kg of body weight of said mammal per day.

50. A method for treating a disease or condition in a mammal, wherein the  $\alpha 7$  nicotinic acetylcholine receptor is implicated comprising administering to a mammal a therapeutically effective amount of compound(s) according to claim 13.
51. The method according to claim 50, wherein the disease or condition is  
5 schizophrenia, or psychosis.
52. The method according to claim 51, wherein said compound(s) is(are) administered rectally, topically, orally, sublingually, or parenterally.
53. The method according to claim 51, wherein said compound(s) is(are) administered from about 0.001 to about 100 mg/kg of body weight of said mammal  
10 per day.
54. The method according to claim 51, wherein said compound(s) is(are) administered from about 0.1 to about 50 mg/kg of body weight of said mammal per day.
55. A method for treating a disease or condition in a mammal, wherein the  $\alpha 7$   
15 nicotinic acetylcholine receptor is implicated comprising administering to a mammal a therapeutically effective amount of compound(s) according to claim 14.
56. The method according to claim 55, wherein the disease or condition is schizophrenia, or psychosis.
57. The method according to claim 56, wherein said compound(s) is(are)  
20 administered rectally, topically, orally, sublingually, or parenterally.
58. The method according to claim 56, wherein said compound(s) is(are) administered from about 0.001 to about 100 mg/kg of body weight of said mammal per day.
59. The method according to claim 56, wherein said compound(s) is(are)  
25 administered from about 0.1 to about 50 mg/kg of body weight of said mammal per day.
60. A method for treating a disease or condition in a mammal, wherein the  $\alpha 7$  nicotinic acetylcholine receptor is implicated comprising administering to a mammal a therapeutically effective amount of compound(s) according to claim 15.
- 30 61. The method according to claim 60, wherein the disease or condition is schizophrenia, or psychosis.
62. The method according to claim 61, wherein said compound(s) is(are) administered rectally, topically, orally, sublingually, or parenterally.

63. The method according to claim 61, wherein said compound(s) is(are) administered from about 0.001 to about 100 mg/kg of body weight of said mammal per day.
64. The method according to claim 61, wherein said compound(s) is(are)  
5 administered from about 0.1 to about 50 mg/kg of body weight of said mammal per day.
65. A method for treating a disease or condition in a mammal, wherein the  $\alpha 7$  nicotinic acetylcholine receptor is implicated comprising administering to a mammal a therapeutically effective amount of compound(s) according to claim 16.
- 10 66. The method according to claim 65, wherein the disease or condition is schizophrenia, or psychosis.
67. The method according to claim 66, wherein said compound(s) is(are) administered rectally, topically, orally, sublingually, or parenterally.
68. The method according to claim 66, wherein said compound(s) is(are)  
15 administered from about 0.001 to about 100 mg/kg of body weight of said mammal per day.
69. The method according to claim 66, wherein said compound(s) is(are) administered from about 0.1 to about 50 mg/kg of body weight of said mammal per day.
- 20 70. A method for treating a disease or condition in a mammal, wherein the  $\alpha 7$  nicotinic acetylcholine receptor is implicated comprising administering to a mammal a therapeutically effective amount of compound(s) according to claim 17.
71. The method according to claim 70, wherein the disease or condition is schizophrenia, or psychosis.
- 25 72. The method according to claim 71, wherein said compound(s) is(are) administered rectally, topically, orally, sublingually, or parenterally.
73. The method according to claim 71, wherein said compound(s) is(are) administered from about 0.001 to about 100 mg/kg of body weight of said mammal per day.
- 30 74. The method according to claim 71, wherein said compound(s) is(are) administered from about 0.1 to about 50 mg/kg of body weight of said mammal per day.

75. A method for treating a disease or condition in a mammal, wherein the  $\alpha 7$  nicotinic acetylcholine receptor is implicated comprising administering to a mammal a therapeutically effective amount of compound(s) according to claim 18.

76. The method according to claim 75, wherein the disease or condition is  
5 schizophrenia, or psychosis.

77. The method according to claim 76, wherein said compound(s) is(are) administered rectally, topically, orally, sublingually, or parenterally.

78. The method according to claim 76, wherein said compound(s) is(are) administered from about 0.001 to about 100 mg/kg of body weight of said mammal  
10 per day.

79. The method according to claim 76, wherein said compound(s) is(are) administered from about 0.1 to about 50 mg/kg of body weight of said mammal per day.

80. A method for treating a disease or condition in a mammal, wherein the  $\alpha 7$  nicotinic acetylcholine receptor is implicated comprising administering to a mammal a  
15 therapeutically effective amount of compound(s) according to claim 19.

81. The method according to claim 80, wherein the disease or condition is schizophrenia, or psychosis.

82. The method according to claim 81, wherein said compound(s) is(are)  
20 administered rectally, topically, orally, sublingually, or parenterally.

83. The method according to claim 81, wherein said compound(s) is(are) administered from about 0.001 to about 100 mg/kg of body weight of said mammal per day.

84. The method according to claim 81, wherein said compound(s) is(are)  
25 administered from about 0.1 to about 50 mg/kg of body weight of said mammal per day.

85. A pharmaceutical composition comprising compound(s) according to claim 1, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.

30 86. A pharmaceutical composition comprising compound(s) according to claim 5, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.

87. A pharmaceutical composition comprising compound(s) according to claim 6, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.
88. A pharmaceutical composition comprising compound(s) according to claim 10, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.
89. A pharmaceutical composition comprising compound(s) according to claim 11, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.
90. A pharmaceutical composition comprising compound(s) according to claim 12, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.
91. A pharmaceutical composition comprising compound(s) according to claim 13, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.
92. A pharmaceutical composition comprising compound(s) according to claim 14, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.
93. A pharmaceutical composition comprising compound(s) according to claim 15, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.
94. A pharmaceutical composition comprising compound(s) according to claim 16, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.
95. A pharmaceutical composition comprising compound(s) according to claim 17, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.
96. A pharmaceutical composition comprising compound(s) according to claim 18, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.
97. A pharmaceutical composition comprising compound(s) according to claim 19, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.

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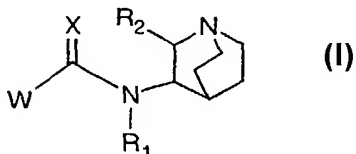
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ance Notes on Codes and Abbreviations" appearing at the begin-  
ning of each regular issue of the PCT Gazette.*

(54) Title: QUINUCLIDINE-SUBSTITUTED HETEROARYL MOIETIES FOR TREATMENT OF DISEASE (NICOTINIC)  
ACETYLCHOLINE RECEPTOR LIGANDS



(57) Abstract: The invention provides compound of Formula (I): These compounds may be in the form of pharmaceutical salts or compositions, and racemic mixtures or pure enantiomers thereof. The compounds of Formula (I) are useful in pharmaceuticals in which  $\alpha 7$  is known to be involved.

# INTERNATIONAL SEARCH REPORT

International Application No

PCT/US 01/22597

## A. CLASSIFICATION OF SUBJECT MATTER

IPC 7 C07D453/02 C07D487/08 A61K31/439 A61P25/00 C07D409/12  
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According to International Patent Classification (IPC) or to both national classification and IPC

## B. FIELDS SEARCHED

Minimum documentation searched (classification system followed by classification symbols)

IPC 7 C07D

Documentation searched other than minimum documentation to the extent that such documents are included in the fields searched

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EPO-Internal, WPI Data, PAJ, CHEM ABS Data

## C. DOCUMENTS CONSIDERED TO BE RELEVANT

Category *	Citation of document, with indication, where appropriate, of the relevant passages	Relevant to claim No.
E	WO 01 60821 A (ASTRAZENECA AB ;PHILLIPS EIFION (US); SCHMIESING RICHARD (US)) 23 August 2001 (2001-08-23) * overlap of chemical formula * the whole document ---	1-97
P,X	WO 01 36417 A (ASTRAZENECA AB ;MULLEN GEORGE (US); PHILLIPS EIFION (US); BALESTRA) 25 May 2001 (2001-05-25) * see claim 1, A = II, definition of R3 * the whole document --- -/--	1-97

☒ Further documents are listed in the continuation of box C.

☒ Patent family members are listed in annex.

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## INTERNATIONAL SEARCH REPORT

International Application No

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## C.(Continuation) DOCUMENTS CONSIDERED TO BE RELEVANT

Category *	Citation of document, with indication, where appropriate, of the relevant passages	Relevant to claim No.
X	PATENT ABSTRACTS OF JAPAN vol. 017, no. 025 (C-1017), 18 January 1993 (1993-01-18) & JP 04 247081 A (TAKEDA CHEM IND LTD), 3 September 1992 (1992-09-03) * overlap of chemical formula * abstract	1-97
X	& JP 04 247081 A (TAKEDA CHEM. IND LTD) 3 September 1992 (1992-09-03) * see page 717, examples *	1-97
X	EP 0 378 111 A (ZAMBON SPA) 18 July 1990 (1990-07-18) * see claim 1, R2 = D, overlap of chemical formula * the whole document	1-97
Y	EP 0 327 335 A (ROBINS CO INC A H) 9 August 1989 (1989-08-09) * see claim 1, definition of Ar * the whole document	1-97
Y	WO 98 54189 A (NIELSEN ELSEBET OESTERGAARD ;NEUROSEARCH AS (DK); OLSEN GUNNAR M ( ) 3 December 1998 (1998-12-03) * see p.2, lines 17 - 33 * the whole document	1-97
Y	WO 97 30998 A (ASTRA AB ;MACOR JOHN (US); WU EDWIN (US)) 28 August 1997 (1997-08-28) * see claim 1, A = IV * the whole document	1-97
Y	HOLLADAY ET AL: "Neuronal Nicotinic Acetylcholine Receptros as Targets for Drug Discovery" JOURNAL OF MEDICINAL CHEMISTRY, AMERICAN CHEMICAL SOCIETY. WASHINGTON, US, vol. 40, no. 26, 19 December 1997 (1997-12-19), pages 4169-4194, XP002124847 ISSN: 0022-2623 * see P.4184, molecular modelling, compare p.4179, ex. 18 and p.4180, ex. 60, 61 * the whole document	1-97
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## INTERNATIONAL SEARCH REPORT

International Application No

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## C.(Continuation) DOCUMENTS CONSIDERED TO BE RELEVANT

Category *	Citation of document, with indication, where appropriate, of the relevant passages	Relevant to claim No.
Y	<p>BANNON A W ET AL: "BROAD-SPECTRUM, NON-OPIOID ANALGESIC ACTIVITY BY SELECTIVE MODULATION OF NEURONAL NICOTINIC ACETYLCHOLINE RECEPTORS" SCIENCE, AMERICAN ASSOCIATION FOR THE ADVANCEMENT OF SCIENCE,, US, vol. 279, no. 5347, 1998, pages 77-81, XP000906771 ISSN: 0036-8075 * see p.77, cpds. fig 1 * -----</p>	1-97

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Information on patent family members

International Application No

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